

#### REASSESSMENT REPORT



### GULF STATES CREOSOTING COMPANY FLOWOOD, RANKIN COUNTY, MISSISSIPPI

U.S. EPA ID No. MSN000407423

Revision 0

#### Prepared for:

U.S. ENVIRONMENTAL PROTECTION AGENCY
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#### 1.0 INTRODUCTION

The U.S. Environmental Protection Agency (EPA) has tasked the T N & Associates, Inc., (TN&A) Superfund Technical Assessment and Response Team (START) to perform a Reassessment under Contract Number (No.) EP-W-05-053 at Gulf States Creosoting Company (Gulf States), EPA Identification (ID) No. MSN000407423, located in Flowood, Rankin County, Mississippi. Reassessments are conducted to evaluate a site's current Hazard Ranking System (HRS) status, document the information contained within the site files, update target information, generate a new site score, and summarize all information in a report submitted to EPA. This Reassessment Report evaluates Gulf States based on site files provided by EPA.

#### 2.0 SITE BACKGROUND

This section describes the site including its environmental, geologic and hydrogeologic setting, historical operations, waste disposal practices, regulatory history, and previous investigations.

#### 2.1 Site Description

The former 141-acre Gulf States Creosoting Company (Gulf States) is located at 1625 Flowood Drive (Mississippi Highway 468), in Flowood, Rankin County, Mississippi (see Appendix A, Figure 1) (Refs. 1, 2). The geographic coordinates at the facility are latitude 32° 18′ 36″ North and longitude 90° 08′ 33″ West (Refs. 2; 3). The facility is currently operating as a horse farm and as ConSteel Co., Inc. (ConSteel) (Refs. 1, p. 2; 4, pp. 2, 3). The horse farm has been operating on the property for at least nine years (Ref. 5, p. 1). The southernmost building on the property is owned by CRC Properties, LLC. (Ref. 4, pp. 2, 5). The northernmost structures including a metal barn and shed are used to house tractors and helicopters for a local news station and are part of the horse farm (Ref. 1, p. 2).

The facility is bound by railroad tracks to the north and east, an adjacent business to the south, and marshland/tributary of the Pearl River to the west. The oxbow lakes, also known as the Creosote Slough, on the western side of the facility are bounded by a large levee, which is fenced and locked (Ref. 5, p. 1). The Creosote Slough can be accessed without going through the fenced area by way of the Pearl River. The Gulf States facility is bound by natural barriers including trees and marshland and is not fenced (Ref. 1, p. 2). The facility is located within a mixed industrial, commercial, and residential area (see Appendix A, Figure 2) (Ref. 2).

ConSteel is a steel fabricator and erector that constructs concrete forms and concrete reinforcement accessories, as well as steel processing and fabrication equipment (Ref. 5, p. 1). File material indicates that ConSteel intends to construct an additional T-shaped office and production facility building on the southwestern portion of the property (Ref. 6, p. 6).

The climate of Rankin County is characterized by long, hot summers and mild winters (Ref. 1, p. 2). Moist tropical air from the Gulf of Mexico has a moderating influence on maximum temperatures in summer (Ref. 7, p. 1). Normal annual total precipitation for the area is approximately 55 inches, and the mean annual lake evaporation is 44 inches, yielding a net annual precipitation of 10 inches (Ref. 8). The 2-year, 24-hour rainfall is 4.5 inches (Ref. 9).

#### 2.2 Geology and Hydrogeology

The facility lies within the Jackson Prairie Belt of the East Gulf Coastal Plain physiographic province (Refs. 7, p. 2; 10, p. 269; 11, p. 23). The property is located along the western border of Rankin County, northeast of Jackson, and east of the Pearl River. The topography of the surrounding area ranges from gently rolling to steep with elevations ranging from 612 feet above mean sea level (amsl) to 270 feet amsl and the property is located at approximately 265 feet amsl (Ref. 2). The property is underlain in descending stratigraphic order by alluvial soils, the Claiborne Group, and the Wilcox Formation (Ref. 13).

The property is directly underlain by alluvial soils of the Cascilla-Arkabutla group, which are classified as nearly level, well-drained to somewhat poorly-drained, silty soils occurring along the flood plains of the Pearl River and its tributaries (Ref. 7, p. 8). The slope of these soils typically ranges from 0 percent (%) to 2% (Ref. 7, p. 8). The nearly linear flood plain surface in the vicinity of the property is irregularly broken by old river runs, natural levees, sloughs, chutes, and scarpes (Ref. 7, p. 8). The average thickness of the alluvium is approximately 40 feet (Ref. 12, p. 32).

The Claiborne Group consists of, in descending stratigraphic order, the Cockfield Formation, the Cook Mountain Formation, the Kosciusko Formation, the Zilpha and Winona Formations, and the Tallahatta Formation (Ref. 13). The Cockfield Formation consists of irregularly bedded laminated lignitic clay, sand, and lignite that is slightly glauconitic (Ref. 13). Based on drillers' logs, the top of the Cockfield is located at approximately 40 feet below land surface (bls) and is approximately 130 feet thick in the vicinity of the property (Ref. 14). The Cook Mountain Formation underlies the Cockfield and consists of marl, limestone, glauconitic sand, and chocholate colored clay (Ref. 13). The Kosciusko Formation, also known as the Sparta Sand, consists of irregularly bedded sand containing clay and small amounts of

quartzite. The Sparta is approximately 300 feet thick in the vicinity of the property (Ref. 11, p. 15). The Zilpha and Winona Formations underlie the Sparta Sand and consist of chocolate colored clay containing glauconitic sand and a highly glauconitic clayey sand, respectively (Ref. 13). The Zilpha ranges in thickness from 200 feet on the Jackson Dome to 420 feet in the southwestern portion of Rankin County (Ref. 11, p. 190). The Winona ranges from 10 to 15 feet thick over the Jackson Dome to approximately 65 feet thick in other portions of the county (Ref. 11, p. 190). The Tallahatta Formation underlies the Zilpha and Winona Formations and consists of predominantly glauconitic claystone and clay with lenses of sand and some sandstone (Ref. 13). The Wilcox Group underlies the Claiborne Group and consists of irregularly bedded fine to coarse sand, more or less lignitic clay, and lignite (Ref. 13). The Wilcox ranges in thickness from approximately 1,100 feet to 1,300 feet over the Jackson Dome and attains a maximum thickness of 2,830 feet in Rankin County (Ref. 11, p. 188).

Three aquifers are available for moderate to large groundwater supplies in Rankin County. The aquifers are, in descending stratigraphic order, the Cockfield Formation, the Sparta Sand, and the Wilcox Group (Ref. 10, p. 274). All of the aquifers are part of the Eocene aquifer system in Mississippi and extend to the west, southwest, and south, and contain freshwater in approximately 50% of the State (Ref. 10, p. 274). All of the aquifers are regional in extent, and all except the Cockfield and lower Wilcox aquifers merge northward into a single aquifer south of Memphis, Tennessee (Ref. 10, p. 274). The formations dip southwest at approximately 15 to 25 feet per mile toward the Mississippi Embayment and the Mississippi River, and the groundwater flow generally follows this regional trend (Ref. 12, p. 4). Within the geologic column, the water-bearing sand beds are interbedded with shale of both marine and continental origin, fossiliferous limestone, and calcareous sandstone (Ref. 12, p. 4). Strata that were deposited by marine origin generally consist of clay and they form aquicludes (confining layers), between the water-bearing sands. These aquicludes are widespread and more uniform in thickness than the aquifers (Ref. 12, p. 4).

The Cockfield Formation is the source of more than half of the municipal water supply in the area, mainly because it is the shallowest of the aquifers (Ref. 12, pp. 1, 32). The Cockfield is an unconfined aquifer located at approximately 40 feet bls in the vicinity of the property, and ranges in thickness from 80 to 140 feet (Refs. 12, p. 32; 14). Municipal water supplies for several small towns are obtained from the Cockfield, some wells yield as much as 500 gallons per minute (gpm) (Ref. 12, p. 1). Based on lithology, the hydraulic conductivity of the Cockfield is approximately 10<sup>-2</sup> centimeters per second (cm/s) (Ref. 15, p. 29). The Cook Mountain underlies the Cockfield and consists of marl, limestone, glauconitic sand, and chocolate colored clay (Ref. 13). The Cook Mountain was deposited in a marine environment, exhibits a high clay content, and serves as a confining layer between the overlying Cockfield aquifer and the

underlying Sparta Sand (Ref. 12, p. 4). Based on lithology, the Cook Mountain exhibits a hydraulic conductivity of approximately 10<sup>-9</sup> cm/s (Ref. 15, p. 29).

The Sparta Sand underlies the Cook Mountain Formation in the vicinity of the property and is approximately 300 feet thick. The Sparta Sand is the most intensively developed aquifer in the vicinity of the property and based on lithology, exhibits a hydraulic conductivity of approximately  $10^{-2}$  cm/s (Refs. 12, pp. 15, 32; 15, p. 29).

The Zilpha and Winona Formations underlie the Sparta Sand and consist of chocolate colored clay containing glauconitic sand and a highly glauconitic clayey sand, respectively (Ref. 13). Due to the high clay content and marine origin of these formations, they serve as the lower confining unit for the Sparta Sand and the upper confining unit for the underlying Wilcox Group (Ref. 12, p. 15). Based on lithology, the Zilpha and Winona Formations exhibit a hydraulic conductivity of approximately 10<sup>-9</sup> cm/s (Ref. 15, p. 29). These formations range in thickness from 420 to 2,600 feet (Ref. 12, p. 15).

The Wilcox Group contains a large reserve of soft water that has been tapped by only a few small-supply wells. The water in this aquifer is more highly mineralized and is warmer than that found in the overlying aquifers (Ref. 12, p. 1). The water is of good quality in Madison and northern Rankin Counties; however, the quality deteriorates down the dip in Hinds County (Ref. 12, p. 1). In counties to the northeast, the results of several pumping tests indicate that the sands in the Wilcox are probably as permeable as the Sparta Sand (Ref. 12, p. 15).

#### 2.3 Ownership and Operations

Gulf States owned the property as early as 1929 and operated a wood treating facility at the location until the mid 1950s (Refs. 1, p. 2; 5, p. 1). During Gulf States operations, railroad cross ties were treated at the facility with coal-tar creosote and transported on and off site using railroad box cars (Ref. 5, p. 1). Coal-tar creosote is a wood preservative used to treat railroad ties, telephone poles, marine pilings, and fence posts (Ref. 16, pp. 1, 2). Contaminants present in coal-tar creosote include polycyclic aromatic hydrocarbons (PAHs), phenol, and cresols (Ref. 16, p. 1). Coal-tar creosote is usually a heavy, oily liquid that is typically amber to brown in color (Ref. 16, p. 2). The creosote found at hazardous waste sites is most often a black, heavy liquid with a sharp smoky odor and burning taste.

In 1958, American Creosoting Corporation purchased portions of the property. In June 1959, W. G. Avery Body Company obtained portions of the property and operated a body shop. ConSteel purchased

the property in 1994 (Ref. 6, p. 2). ConSteel currently owns the 8-acre property on which the former operations area of the Gulf States facility lies (Ref. 4, pp. 2, 3). ConSteel appears to be operating on the portion of the property that it owns. Avery Lead Track, LLC., currently owns two additional parcels of the property, totaling approximately 120 acres (Ref. 4, pp. 2, 4, 6, 7).

#### 2.4 Regulatory and Release History

In July 1993, a Phase I Environmental Assessment (Phase I) of the property included a review of the Mississippi Department of Environmental Quality (MDEQ) Underground Storage Tank (UST) division records (Ref. 1, p. 3). No past UST usage on the property was found during the review. No other regulatory or permitting information has been identified for the facility.

#### 2.5 Previous Investigations

In June 1993, BCM Engineers, Inc. (BCM) conducted a Phase I of the former Gulf States property on behalf of Trustmark National Bank (Ref. 1, p. 3). The objective of the Phase I was to identify adverse environmental conditions, suspect activities, and potential hazardous wastes or materials on or in the vicinity of the subject property. The Phase I included the collection of soil samples from soil borings ranging in depth from 0 to 8 feet below ground surface (bgs). The seven soil borings were concentrated in the suspected location of the former creosote operations. Of the seven borings, five samples were collected and analyzed for semi-volatile organic compounds (SVOCs). Constituents detected in the soil samples were naphthalene, 2-methylnaphthalene, dibenzofuran, fluorene, phenanthrene, and fluoranthene. The concentrations ranged from below the detection limit to 604 milligrams per kilogram (mg/kg) (Ref. 1, p. 3).

In August 1993, BCM conducted a Phase II of the former Gulf States property (Ref. 1, p. 3). The objective of the Phase II was to delineate the extent of the creosote-contaminated soil identified during the Phase I investigation. During the Phase II, several soil borings were advanced and five soil samples were collected for SVOC laboratory analysis (Ref. 1, p. 4). The soil borings ranged in depth from 0 to 8 feet bgs. Sample results indicated the presence of 2-methylnaphthalene, naphthalene, dibenzofuran, phenanthrene, fluoranthene, pyrene, and chrysene. Constituents ranged from below the detection limit to 1,057 mg/kg. The data were compared to health-based criteria for exposure via groundwater ingestion. Based on such a comparison, the recommendation for no further action was presented because the soil contaminants were below the target cleanup levels (Ref. 1, p. 4).

In April 2003, Weston Solutions, Inc. (Weston) conducted a Preliminary Assessment/Site Inspection (PA/SI) at the site for EPA (Ref. 1). All sampling was conducted by the EPA Science and Ecosystem Support Division (SESD) (Ref. 17). SESD collected surface soil samples, subsurface soil samples, groundwater samples, and sediment samples on or near the Gulf States property (see Appendix A, Figures 3 and 4) (Refs. 1, p. 4; 17). Sampling specifics are detailed in Section 2.5.1.

In May 2006, EPA SESD and MDEQ collected fish, sediment, and surface water samples within the Creosote Slough located on the western side of the property and analyzed them for PAHs. The study area was comprised of a reference station (CS-01) and three other stations (CS-02, CS-03, and CS-04) that covered the area from the railroad tracks on the north end of the property to the terminus of the slough at the Pearl River (Ref. 5, p. 1).

PAH concentrations in surface water were compared to the Ambient Water Quality Criteria for Human Health (AWQC-HH) (Ref. 5, p. 2). PAH concentrations in sediments were compared to EPA Region 9 Preliminary Remediation Goals (PRGs) for industrial soil (Ref. 5. p. 3). In all fish tissue samples, an EPA reference titled *Guidance for Assessing Chemical Contaminant Data for use in Fish Advisories, Vol. 1* (US EPA 2000), was used in assessment of the results. In summary, levels of PAHs detected in surface water, sediment, and fish tissue samples during the May 2006 sampling event were found to be of no current concern for human exposure (Ref. 5, p. 3).

In September 2006, Earth Consulting Group, Inc. (EarthCon) performed a Limited Soil Assessment of the western portion of the ConSteel property located at 1625 Flowood Drive. The assessment included the installation of soil borings and selected soil sampling and analysis to assess the area for the presence of creosote compounds related to Gulf States prior to the purchase of the property by ConSteel (Ref. 6, p. 1).

An investigation related to a pending commercial transaction for ConSteel discovered a small depression in the northwest corner of the property that contained remnants of a concrete culvert (see Appendix A, Figure 5) (Ref. 6, p. 2). Borings B-1 thorough B-12, B-17 through B-20, B-26, and B-27 were installed with a Geoprobe<sup>®</sup> in a general radial pattern away from this area to attempt to delineate the lateral extent of the creosote impacts associated with this feature (Ref. 6, p. 2). Other areas of potential concern on the property included a possible buried drain line leading from the west side of the existing ConSteel production building to the concrete culvert remnant in the northwest corner of the property. Soil borings B-13 through B-16 were installed along the suspected route of this potential drain line including areas of stressed vegetation observed along the route (Ref. 6, p. 2). Soil borings B-21 through B-25 were installed within the footprint of a planned new ConSteel production and office facility on the southwest portion of

the property. These borings were installed to assess possible residual creosote impacts to the subsurface beneath the area where workers would be present during planned future operations (Ref. 6, pp. 2, 3). A total of 27 soil borings were installed to a depth of approximately 8 feet bgs (Ref. 6, p. 3).

According to EarthCon, approximately 15,000 square feet of the northwest corner of the property appeared to be impacted with creosote odors and/or soil staining in the shallow subsurface (Ref. 6, p. 5). Free creosote product was observed in soil borings at two locations on the site (Ref. 6, p. 5). These areas included a depression near the northwest corner of the subject property, and an area containing buried wood debris in the west-central portion of the subject property (Ref. 6, pp. 5, 6). Borings installed in the shallow depression in the northwest corner of the property were observed to contain degraded, soft soils and liquid creosote product in the upper few feet, with heavy creosote odors and stained soil to the termination depth of the borings at approximately 8 feet bgs (Ref. 6, p. 5). Creosote-soaked wood debris was encountered in several boring locations at a depth of approximately 3.5 to 4.5 feet bgs (Ref. 6, p. 6). The thickness of the buried wood debris in this area was not defined. Laboratory analytical results performed by Environmental Science Corporation (ESC) of soil samples collected from each of the two identified creosote locations on site contained SVOCs (see Appendix B, Table 1). Contaminants including benzo(a)anthracene, chrysene, and naphthalene were detected at concentrations exceeding Region 9 PRGs for industrial soil (Ref. 6, pp. 4, 5, 8).

On September 20, 2007, EPA, SESD, and MDEQ conducted a Removal Site Evaluation (RSE) at the site to ascertain the existence of a buried pipe system and determine whether the system could be contributing to contamination found in oxbow lakes (Creosote Slough) just west of the site (Refs. 18, p. 1; 19). Although an underground pipe network was identified and traced from the site to the marsh, there was no evidence of the pipe being a conduit for creosote from the former treatment area onto the marsh. The RSE revealed no trace of creosote inside the pipe (Ref. 19).

#### 2.5.1 2003 PA/SI

#### Surface and Subsurface Soil Samples

SESD collected 24 surface soil samples and 21 subsurface soil samples at Gulf States in 2003 (Refs. 1, pp. 6–9; 17). One background surface soil sample (GS-01-SS) and one background subsurface soil sample (GS-01-SB) were collected northeast of the facility from Jackson Preparatory School at 3100 Lakeland Drive. SESD collected the surface soil samples from 0 to 6 inches bgs, and subsurface soil samples were collected from 2 to 3 feet bgs with the exception of two locations, which were collected between 6 and 12 inches (GS-10-SB and GS-14-SB) (Ref. 17, p. 2). For purposes of this reassessment

and adherence to the HRS, all samples collected from 0 to 2 feet bgs are considered surface soil samples. Samples collected at depths greater than 2 feet bgs are considered subsurface soil samples.

On site surface and subsurface soil samples were collected over the entire property (see Appendix A, Figure 4). Surface soil analytical results revealed HRS-elevated concentrations (three times background) of several inorganic contaminants including antimony, barium, beryllium, cadmium, copper, lead, manganese, magnesium, nickel, thallium, and zinc (see Appendix B, Table 2). No inorganic contaminants were detected at levels above the Region 9 PRG for industrial soil. Organic compounds were detected at HRS-elevated concentrations in the surface soil. These compounds include benzo(b)fluoranthene, acenaphthylene. anthracene. benzo(a)anthracene. benzo(ghi)pervlene. benzo(k)fluoranthene, benzo(a)pyrene, carbazole, chrysene, dibenzo(a,h)anthracene, fluoranthene, indeno(1,2,3-cd)pyrene, naphthalene, phenanthrene, pyrene, 4,4'-DDE, 4,4'-DDT, endrin, endrin aldehyde, endrin ketone, methoxchlor, and methyl ethyl ketone (see Appendix B, Table 3) (Ref. 17). Several organic contaminants were above EPA Region 9 PRGs for industrial soil including benzo(a)anthracene, benzo(b)fluoranthene, benzo(k)fluoranthene, benzo(a)pyrene, dibenzo(a,h)anthracene, and indeno(1,2,3-cd)pyrene.

Elevated contaminants detected in the subsurface soil include manganese, selenium, benzo(b)fluoranthene, benzo(k)fluoranthene, bis(2-ethylhexyl)phthalate, indeno(1,2,3-cd)pyrene, acetone, 4,4'-DDE, and Aroclor 1260 (see Appendix B, Tables 4 and 5) (Ref. 17).

#### **Groundwater Samples**

SESD collected five groundwater samples from temporary monitoring wells installed at Gulf States (see Appendix A, Figure 4) (Ref. 1, pp. 10, 19). The background sample was collected from the Jackson Preparatory School at 3100 Lakeland Drive located northeast of the facility (see Appendix A, Figure 3). On-site well locations are described in the PA/SI (Ref. 1).

Inorganic and organic contaminants were detected at elevated concentrations in the on-site temporary monitoring wells (see Appendix B, Tables 6 and 7). These contaminants include aluminum, barium, beryllium, cobalt, iron, manganese, magnesium, nickel, vanadium, zinc, and methyl ethyl ketone (Ref. 17).

#### Sediment Samples

Eight sediment samples were collected during the PA/SI to document the migration of on-site contaminants into the marsh located west of the Gulf States property, and the Pearl River (see Appendix

A, Figure 4) (Ref. 1, p. 11). The background sample (GS-01-SD) was collected northeast of the site behind Jackson Preparatory School (see Appendix A, Figure 3) (Refs. 1; 17, p. 2). Control samples were collected upstream of the site on the Pearl River (GS-06-SD) and on Prairie Branch (GS-08-SD) to isolate various potential influences (Ref. 17, p. 2). Since a sediment sample was not collected for comparison to control sample GS-08-SD, this sample was not used for this evaluation.

Elevated contaminants detected in sediment samples from the marsh include barium, beryllium, lead, selenium, zinc, benzo(b)fluoranthene, benzo(k)fluoranthene, 2-methylnaphthalene, acenaphthene, acenaphthylene, anthracene, benzo(a)anthracene, benzo(a)pyrene, chrysene, dibenzo(a,h)anthracene, fluoranthene, fluorene, naphthalene, phenanthrene, pyrene, 4,4'-DDE, beta-BHC, acetone, carbon disulfide, and methyl ethyl ketone (see Appendix B, Tables 8 and 9) (Ref. 17). Of these contaminants, lead, zinc, and 4,4'-DDE were above their respective sediment screening values.

No elevated contaminants were detected in the Pearl River sediment samples (see Appendix B, Tables 10 and 11) (Ref. 17).

#### 2.6 Potential Source Areas

The potential source area at Gulf States is contaminated soil. Contaminated soil has been detected throughout the property and comprises approximately 1,441,876 square feet (or 33 acres). Based on analytical results from 2003 and 2006, SVOCs and pesticides including acenaphthylene, anthracene, benzo(a)anthracene, benzo(b)fluoranthene, benzo(ghi)perylene, benzo(k)fluoranthene, benzo(a)pyrene, carbazole, chrysene, dibenzo(a,h)anthracene, fluoranthene, indeno(1,2,3-cd)pyrene, naphthalene, phenanthrene, pyrene, 4,4'-DDE, 4,4'-DDT, endrin, endrin aldehyde, endrin ketone, and methoxychlor were present in the surface soil as a result of Gulf States' operations (Refs. 1, 17, 16, 6). Of these contaminants, benzo(b)fluoranthene, benzo(k)fluoranthene, indeno(1,2,3-cd)pyrene, and 4,4'-DDE have migrated to the subsurface soil at elevated concentrations.

Although numerous metals were detected in surface soil and groundwater samples, the contamination cannot be attributed to Gulf States operations. The metals contamination is likely a contaminant derived from the body shop activities from W.G. Avery Body Company or current site activities from ConSteel.

#### 3.0 PATHWAYS

This section discusses the groundwater migration, surface water migration, soil exposure, and air migration pathways. Additionally, this section discusses the targets associated with each pathway and draws pathway-specific conclusions. Sampling locations and analytical results for samples collected from the specific pathways are also discussed.

#### 3.1 Groundwater Migration Pathway

Numerous metals and one VOC (methyl ethyl ketone) were detected in on-site groundwater samples; however, the groundwater contamination cannot be attributed to Gulf States.

The groundwater migration pathway is of potential concern at Gulf States since many nearby residents obtain their drinking water from groundwater sources. Based on information provided by MDEQ, nearby residents within 4-miles of the site obtain drinking water from the City of Flowood, the City of Pearl, the City of Richland, the City of Jackson, and Cleveland's Trailer Park (Ref. 20). Drinking water is obtained from groundwater wells screened in the Cockfield and Sparta aquifers, with the exception of the City of Jackson, who services a surface water intake. This intake will be discussed further in the surface water pathway.

The City of Flowood, the City of Pearl, and the City of Richland draw their water from the Sparta Aquifer (Ref. 20). Several wells from these water departments are located within 4 miles of the site (Ref. 2). Two City of Flowood wells are located within the 0.5 mile radius of the site; however, these two wells are currently inactive (reasons unknown) and were not considered for HRS scoring purposes in this assessment (Ref. 20). The City of Flowood operates two wells in the vicinity of the property; one well located within the 0.5 to 1 mile radius, and one well within the 1 to 2 mile radius. Each City of Flowood well maintains 1,500 connections (or 3,930 persons) (Refs. 20, 22). The City of Pearl operates three wells located within the 2 to 3 mile radius from the site. Each City of Pearl well maintains 7,351 connections (or 19,259 persons) (Refs. 20, 22). One additional City of Pearl well is located within the 3 to 4 mile radius of the site. The City of Richland operates one well located within the 3 to 4 mile radius of the site. The City of Richland well maintains 2,157 connections (or 5,651 persons) (Refs. 20, 22). All municipal wells in the area are located within designated Wellhead Protection Areas (Ref. 20).

Cleveland's Trailer Park maintains four community wells located within 2 to 3 miles from the site (Ref. 20). These wells are screened in the Cockfield Aquifer. Each of Cleveland's Trailer Park wells maintains 135 connections (or 353 persons) (Refs. 20, 22).

It is unknown whether private wells are present within a 4-mile radius of Gulf States. No private well information was available through the water departments or MDEQ at the time of inquiry for this reassessment. However, Weston noted in September 2003 that no private wells were observed during the site reconnaissance for the PA/SI (Ref. 21).

#### 3.2 Surface Water Migration Pathway

The City of Jackson obtains municipal water from a surface water intake; however, the intake is not located within the 15-mile surface water pathway from the facility (Ref. 20). Several endangered and threatened species are present within the State of Mississippi; however, no file material is available indicating that any of these species are present in the surface water migration pathway of the site (Ref. 23). The property is not located within a flood plain (Ref. 24). There is no evidence of HRS-qualifying wetlands along the surface water migration pathway from the site (Ref. 25).

Sampling of the surface water migration pathway does not indicate HRS elevated levels of site-attributable constituents and no sampling data from the release is available in the file material. Although sediment samples have been collected from the Pearl River, no compounds were detected at HRS-elevated concentrations; therefore, no observed release to the surface water pathway has been documented. Additionally, the target values for the surface water migration pathway are extremely low. Therefore, the surface water migration pathway at Gulf States is of minimal concern.

#### 3.3 Soil Exposure Pathway

The property is surrounded by natural barriers and a well-maintained fence. No residences are present on sources at the site and the number of nearby residents associated with the soil exposure pathway is minimal. The ConSteel facility is active; however, the number of workers is unknown. Therefore, the soil exposure pathway is of minimal concern and was not evaluated for this assessment.

#### 3.4 Air Migration Pathway

No air samples have ever been collected from the Gulf States facility and no releases to the air migration pathway have been documented. Therefore, the air migration pathway at Gulf States is of minimal concern and was not evaluated for this reassessment.

#### 4.0 CONCLUSIONS

The former 141-acre Gulf States Creosoting Company (Gulf States) is located at 1625 Flowood Drive (Mississippi Highway 468), in Flowood, Rankin County, Mississippi. Gulf States owned the property as early as 1929 and operated a wood treating facility at the location until the mid 1950s. During Gulf States operations, railroad cross ties were treated at the facility with coal-tar creosote and transported on and off site using railroad box cars. In 1958, American Creosoting Corporation purchased portions of the property. In June 1959, W. G. Avery Body Company obtained portions of the property and operated a body shop. ConSteel purchased the property in 1994. ConSteel currently owns the 8-acre property on which the former operations area of the Gulf States facility lies. ConSteel appears to be operating on the portion of the property that it owns. Avery Lead Track, LLC., currently owns two additional parcels of the property, totaling approximately 120 acres.

Several investigations have been conducted at the facility including a Phase I in June 1993 by BCM on behalf of Trustmark National Bank. The Phase I was conducted to identify any adverse environmental conditions, suspect activities, and potential hazardous wastes or materials on or in the vicinity of the property. During the Phase I, soil samples were collected from soil borings ranging in depth from 0 to 8 feet below ground surface (bgs). Contamination was identified, and in August 1993, BCM conducted a Phase II to delineate the extent of the creosote-contaminated soil identified during the Phase I investigation. During the Phase II, several soil borings were advanced and five soil samples were collected for SVOC laboratory analysis. The soil borings ranged in depth from 0 to 8 feet bgs.

In April 2003, Weston conducted a PA/SI at the site for EPA. During the PA/SI, SESD collected surface soil samples, subsurface soil samples, groundwater samples, and sediment samples on or near the Gulf States property. Results of the investigation indicated the presence of several SVOC and PAH in the surface and subsurface soils at the site.

In May 2006, EPA SESD and MDEQ collected fish, sediment, and surface water samples within the Creosote Slough located on the western side of the property for PAHs analysis. Levels of PAHs detected

in surface water, sediment, and fish tissue samples during the sampling event were found to be of no current concern for human exposure.

In September 2006, EarthCon performed a Limited Soil Assessment including the installation of soil borings and selected soil sampling and analysis to assess the area for the presence of creosote compounds related to Gulf States prior to the purchase of the property by ConSteel. These borings were installed to assess possible residual creosote impacts to the subsurface beneath the area where workers would be present during planned future operations. A total of 27 soil borings were installed to a depth of approximately 8 feet bgs. According to EarthCon, approximately 15,000 square feet of the northwest corner of the property appeared to be impacted with creosote odors and/or soil staining in the shallow subsurface. Contaminants including benzo(a)anthracene, chrysene, and naphthalene were detected at concentrations exceeding Region 9 PRGs for industrial soil.

On September 20, 2007, EPA, SESD, and MDEQ conducted a RSE at the site to ascertain the existence of a buried pipe system and determine whether the system could be contributing to contamination found in oxbow lakes (Creosote Slough) just west of the site. Although an underground pipe network was identified and traced from the site to the marsh, there was no evidence of the pipe being a conduit for creosote from the former treatment area onto the marsh. The RSE revealed no trace of creosote inside the pipe.

The groundwater migration pathway is the only pathway of concern at the Gulf States property; however, no observed release to the groundwater pathway has been documented and the number of potential targets associated with the pathway is relatively low. Several inorganic constituents have been identified in soil samples and in groundwater beneath the property; however, metals are not attributable to the operations by Gulf States at the facility. The presence of metals in the groundwater is likely a result of operations by ConSteel or Avery. Both aquifers underlying the property were evaluated as part of this reassessment; however, due to the absence of an observed release to groundwater, all targets associated with the groundwater migration pathway are potential targets.

The surface water pathway at the facility is of minimal concern because no observed release to the surface water pathway has been documented and the number of targets associated with the pathway is minimal.

The soil exposure pathway at the facility is of minimal concern because access to the facility is hindered by natural barriers and a well maintained chain-linked fence. Additionally, the number of workers is unknown, no residences are located on source areas, and the number of nearby potential targets is minimal.

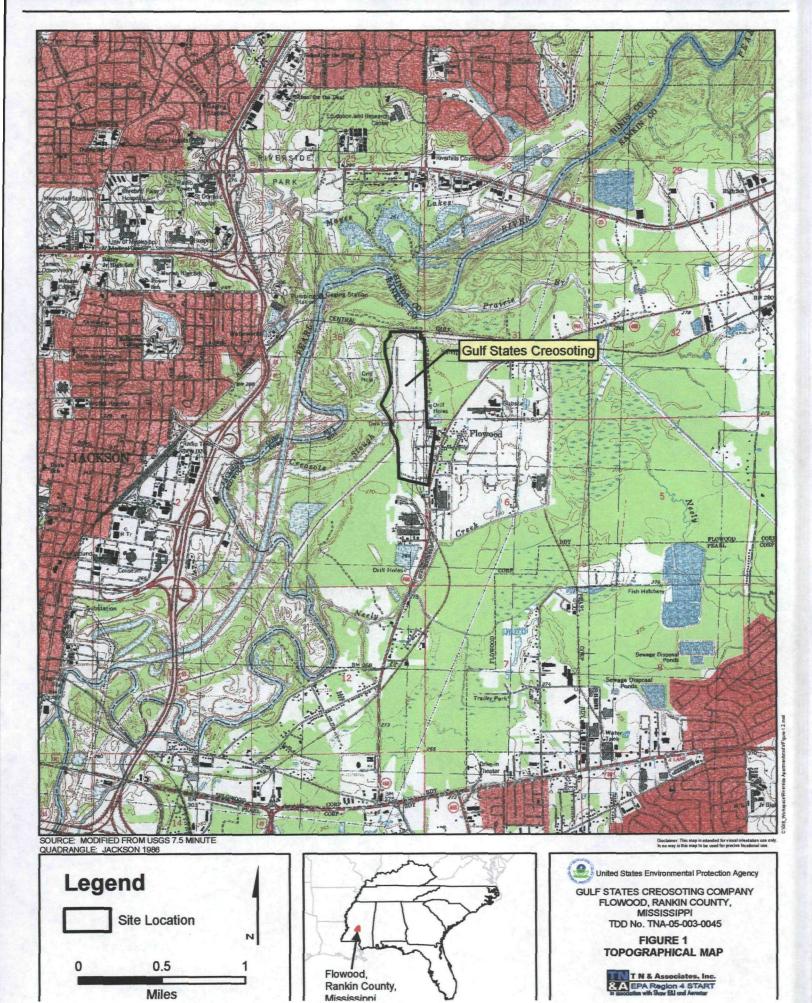
The air migration pathway at the facility is of minimal concern. No air samples have been collected from the facility and no observed release to the air migration pathway has been documented.

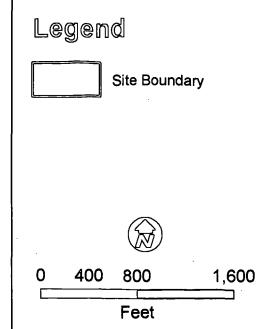
Based on the currently available information, the Gulf States Creosoting site score is 26.8. This score is less than the cutoff value of 28.50 necessary to consider listing a site on the National Priorities List (NPL). EPA will determine the need for further remedial actions at this property.

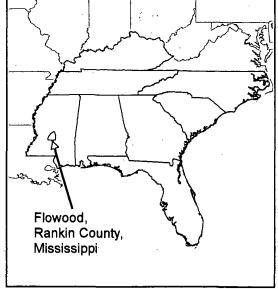
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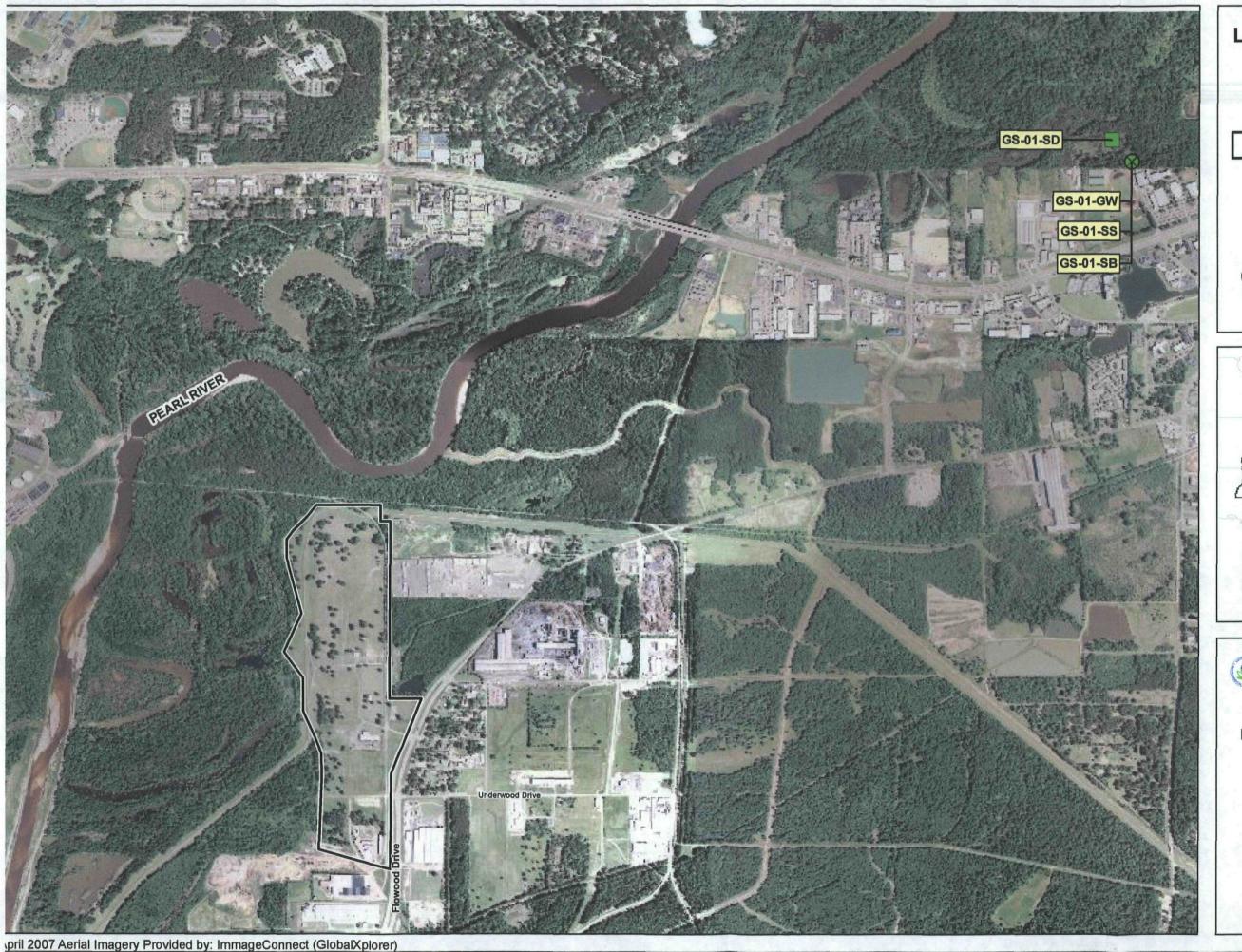


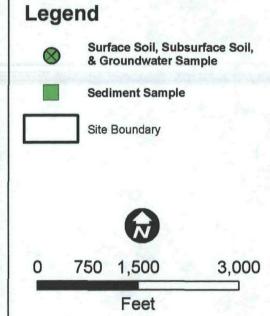
United States Environmental Protection Agency

**GULF STATES CREOSOTING COMPANY** FLOWOOD, RANKIN COUNTY, MISSISSIPPI TDD No. TNA-05-003-0045

FIGURE 2 AERIAL PHOTOGRAPH

T N & Associates, Inc.



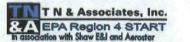


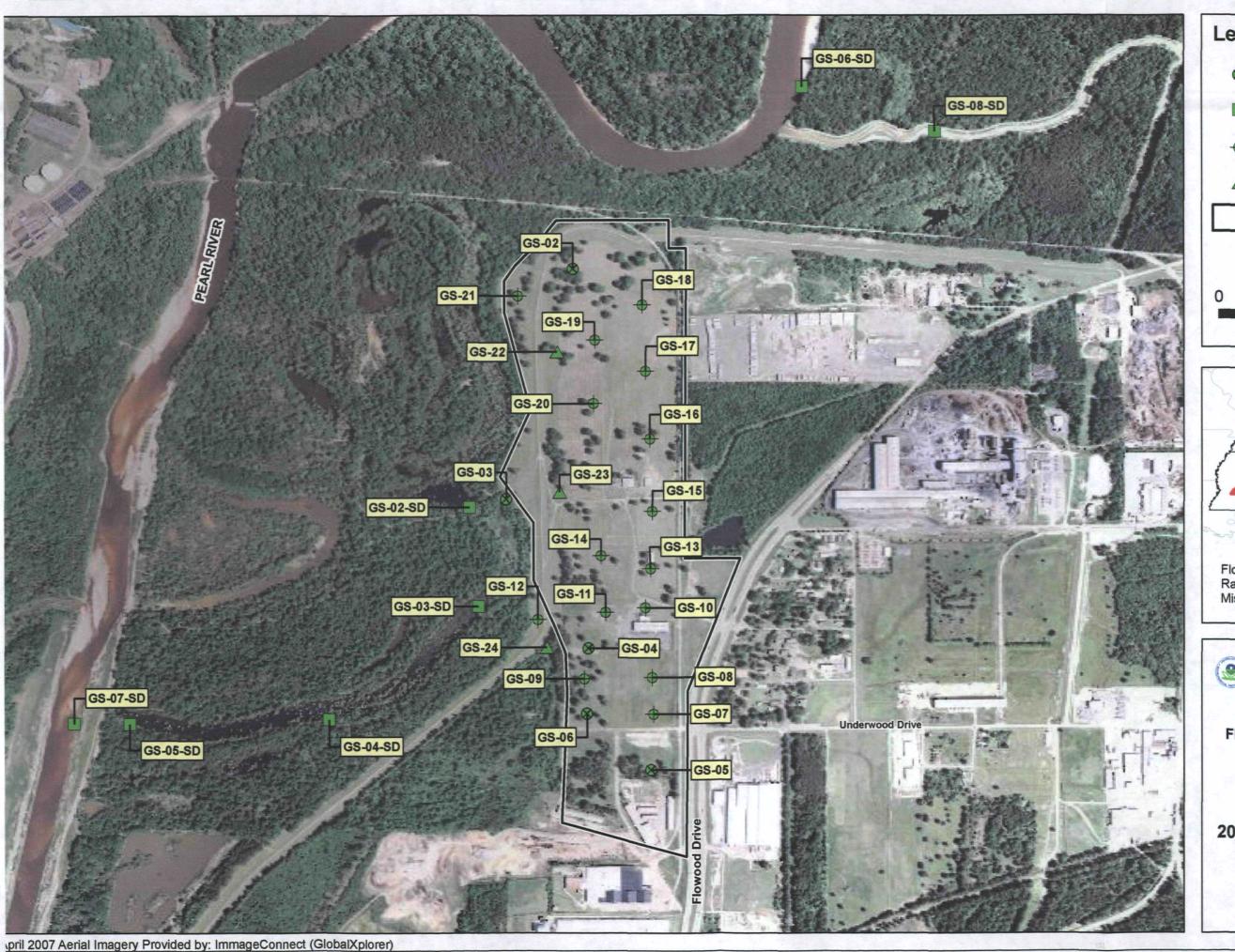


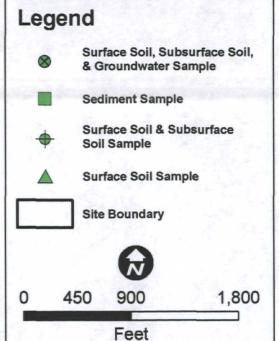


**GULF STATES CREOSOTING COMPANY** FLOWOOD, RANKIN COUNTY, MISSISSIPPI TDD No. TNA-05-003-0045

FIGURE 3 **2003 BACKGROUND SAMPLE LOCATIONS** 





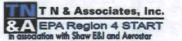




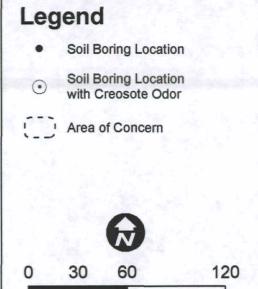


**GULF STATES CREOSOTING COMPANY** FLOWOOD, RANKIN COUNTY, MISSISSIPPI TDD No. TNA-05-003-0045

FIGURE 4 **2003 SAMPLE LOCATIONS** 

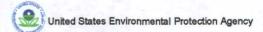






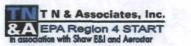
Feet



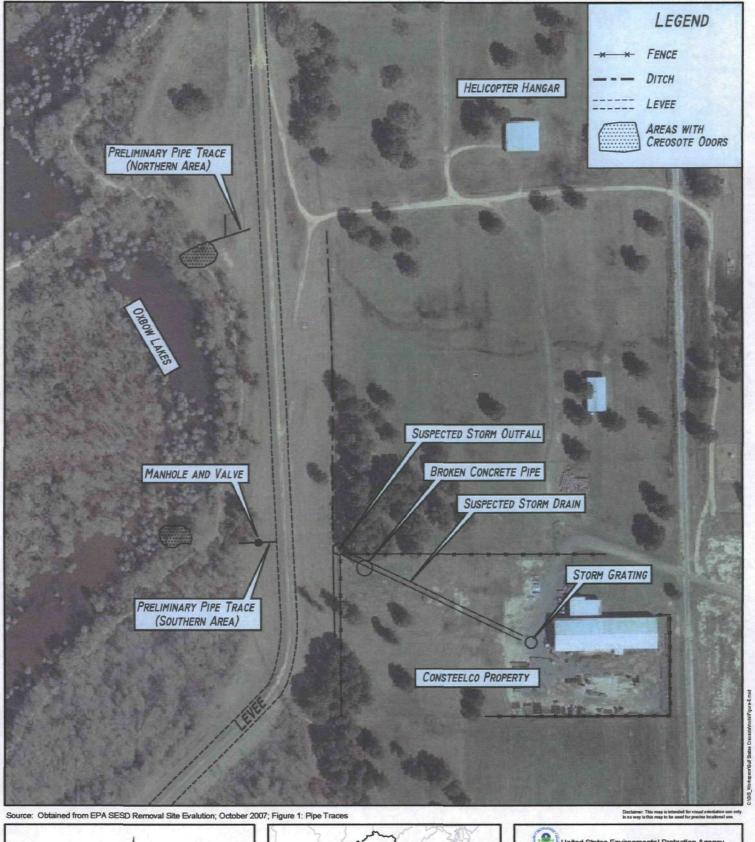


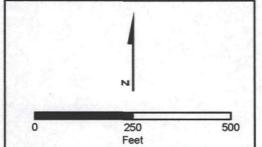
**GULF STATES CREOSOTING COMPANY** FLOWOOD, RANKIN COUNTY, MISSISSIPPI TDD No. TNA-05-003-0045

FIGURE 5 **2006 SAMPLE LOCATIONS** 

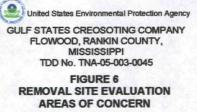


Source: Obtained from Earth Consulting Group, Inc. Limited Soil Assessment; October 2006; Figure 2: 2003 Aerial Photograph-Soil Boring Location Map

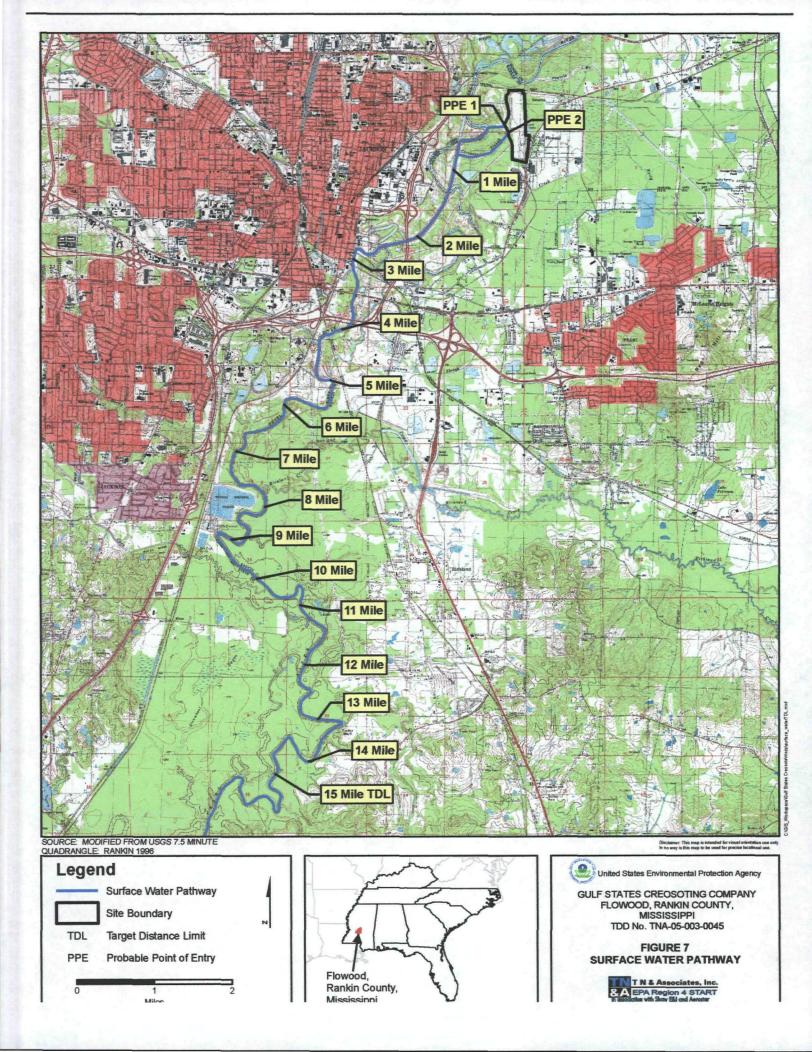












APPENDIX B

Tables

TABLE 2
SUMMARY OF 2003 SURFACE SOIL INORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-02-SS	GS-03-SS	GS-04-SS	GS-05-SS	GS-06-SS	GS-07-SS	GS-08-SS	PRG
ANALTYE	Background				•				1110
Metals, Total (mg/k	g)	_						<del>.</del>	
Aluminum	4700 J	7800 J	8900 J	4300 J	3100 J	6700 J	2800 J	1800 J	100000
Antimony	0.56 UJ	0.64 R	0.64 U	0.63 U	0.57 U	0.65 U	0.58 U	0.5 U	409
Arsenic	6.3	3.3 J	11	2	13	2	3.6	3.4	1.59
Barium	72	120	67	60	37	85	34	21	66600
Beryllium	0.33	0.94	0.94	0.37	0.3	0.28	0.2	0.19	1940
Cadmium	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U	0.05 U	0.09	451
Chromium	8.4	6.1	20	6	7.5	7.7	4.5	4	448
Cobalt	5.8	17	8.1	1.7	6.6	1.4	2	1.5	1920
Copper	6	2 J	6.5 U	3.9 U	3 U	4.9 U	3.5 U	2.9 U	40900
Iron	10000	8400	27000 J	4500 J	15000 J	5600 J	5400 J	4700 J	100000
Lead	11 J	17	1203 完全	18	19	17	11	12	800
Magnesium	520 J	470 J	680	280	180	550	250	180	NL
Manganese	630	2800	550	140	610	170	180	230	19500
Nickel	4.8 U	8.3 U	5.1 U	2.6 U	2 U	2.8 U	1.9 U	1.7 U	20400
Selenium	1.3	0.93 R	1.9 J	0.63 U	0.87 U	0.65 U	0.79 U	0.68 U .	5110
Silver	0.39 R	0.71 R	0.63	0.26 R	0.43	0.29	0.17 U	0.15 U	5110
Thallium	0.73 U	0.81 U	0.83 U	0.81 U	0.73 U	0.84 U	0.75 U	0.65 U	67.5
Vanadium	17	15	46	11	18	13	8.1	7.1	1020
Zinc	29	47	34	27	23	35	24	26	100000

Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SS - Surface soil

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

U - Analyte was not detected at or above the detection limit

# TABLE 2 SUMMARY OF 2003 SURFACE SOIL INORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-09-SS	GS-10-SB*	GS-10-SS	GS-11-SS	GS-12-SS	GS-13-SS	GS-14-SB*	PRG
ANALYTE	Background				·				1 KO
Metals, Total (mg/kg)									
Aluminum	4700 J	2500 J	8600 J	4700 J	4400 J	5500 J	4400 J	4600 J	100000
Antimony	0.56 UJ	1.2 R	0.96 R	. 1R	0.51 UJ	0.58 U	0.6 R	0.56 ÚJ	409
Arsenic	6.3	0.95 R	13	6.8	3.7	4.4	6.3	7.4	. 1.59
Barium	72	25	4301:54.~	120	54	83	58	54	66600
Beryllium	0.33	0.15		0.46	0.44	0.46	0.54	0.48	1940
Cadmium	0.05 U	0.05 U	0.87	0.34	0.19 🏕 🔻	0.05 U	0.16 R	0.05 U	451
Chromium	8.4	4.1	19	23	12	8.9	7.8	7.1	448
Cobalt	5.8	0.95	12	4.8	5.4	7.3	4.2	12	1920
Copper	6	0.87 UJ	್: 40 J 🖏	19 J	5 UJ_	3.4 U	7.2 J	1.5 UJ	40900
Iron	10000	2800	20000	14000	11000	11000 J	12000	14000	100000
Lead	11 J	8.8	ALTERNATION 18	80 🗼 🗼	290 🦀 🎎	25	. 33	27	800
Magnesium	520 J	160 J	2400 J 🧎 📀	780 J	260 J	380	280 J	250 J	NL
Manganese	630	28	1800	500	880	990	940	1000	19500
Nickel	4.8 U	1.5 U	20 至分	8.5 U	4.9 U	3.2 U	5.2 U	3.3 U	20400
Selenium	1.3	0.6 U	1.2	0.98 R	0.77 R	1.1 U	1.2	0.56 U	5110
Silver	0.39 R	0.28 R	0.76	0.49	0.38	0.43	0.33 R	0.42	5110
Thallium	0.73 U	0.77 U	0.76 U	0.7 U	0.66 U	0.75 U	0.69 U	0.73 U	67.5
Vanadium	17	6.6	30	18	16	22	17	23	1020
Zinc	29	19	290	160	<b>840</b>	28	72	20	100000

Notes:

Shaded - Concentration is elevated (3 x background)

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SB - Subsurface soil

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mg/kg - Milligrams per kilogram

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

U - Analyte was not detected at or above the detection limit

# TABLE 2 SUMMARY OF 2003 SURFACE SOIL INORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-14-SS	GS-15-SS	GS-16-SS	GS-17-SS	GS-18-SS	GS-19-SS	GS-20-SS	PRG ·
ANALYTE	Background								
   Metals, Total (mg/k	g)								
Aluminum	4700 J	4800 J	8000 J	7700 J	9000 J	4900 J	8500 J	6700 J	100000
Antimony	0.56 UJ	∷0.69 J¢	0.56 UJ	0.7 UJ	0.83 UJ	1 UJ	0.89 UJ	0.6 UJ	409
Arsenic	6.3	9	6.9	5.5 J	5.5	5.4	6.5	5.8	1.59
Barium	72	65	72	50	120	53	89	410	66600
Beryllium	0.33	0.44	0.71	0.43	0.94	0.35	0.73	1.2	1940
Cadmium	0.05 U	0.11 R	0.1	0.05 U	0.07 R	0.05 U	0.05 U	0.17	451
Chromium	8.4	8.6	11	10	8	9.9	8.3	8.2	448
Cobalt	5.8	7.3	11	4	12	7.6	13	17	1920
Copper	6	5.7 UJ	4.8 UJ	5.3 U	5.5 U	3.1 U	3.8 U	0.95 U	40900
Iron	10000	11000	12000	12000	12000	10000	12000	12000	100000
Lead	11 J	24	26	15 J	20 J	15 J	19 J	24 J	800
Magnesium	520 J	340 J	430 J	510 J	490 J	340 J	530 J	380 J	NL
Manganese	630	870	1800	240	2500	590 .	鸞∄1900	5500	19500
Nickel	4.8 U	4.1 U	6.7 U	3.6 U	9.2 U	3.9 U	6.7 U	8 U	20400
Selenium	1.3	0.63 R	1.1 R	1.3	1.5 J	1.1	1.2	1.3 J	5110
Silver	0.39 R	0.37	0.5	0.39 R	0.51 R	0.29	0.53 R	. 0.9 R	5110
Thallium	0.73 U	0.72 U	0.73 U	0.78 U	0.79 U	0.8 U	0.73 U	· 1.5 《 · ) ※:	67.5
Vanadium	17	21	23	22	23	- 20	22	22	1020
Zinc	29	44	48	33	65	34	52	88	100000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SS - Surface soil

SB - Subsurface soil

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J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

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# TABLE 2 SUMMARY OF 2003 SURFACE SOIL INORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-21-SS	GS-22-SS	GS-23-SS	GS-24-SS	PRG
ANALYTE	Background					
Metals, Total (mg/kg	1)			· · · · · ·		
Aluminum	4700 J	6800 J	4200 J	7500 J	4900 J	100000
Antimony	0.56 UJ	0.96 UJ	0.56 U	0.65 UJ	0.91 U	409
Arsenic	6.3	2.7	4	5	6.5	1.59
Barium	72	88	42	83	120	66600
Beryllium	0.33	0.81	0.43	0.5	0.61	1940
Cadmium	0.05 U	0.05 U	0.05 U	0.1 R	0.16	451
Chromium	8.4	5.6	9.4	12	12	448
Cobalt	5.8	14	5.5	6.8	. 6.8	1920
Copper	6	1.5 U	2.5 U	12 J	8.4 U	40900
Iron	10000	7300	9700 J	13000	17000 J	100000
Lead	11 J	13 J	9.8	24	28	800
Magnesium	520 J	270 J	220	630 J	450	NL
Manganese	630	2200	470	920	890	19500
Nickel	4.8 U	6.7 U	2.2 U	4.8 U	5.3 U	20400
Selenium	1.3	0.8	1.1 U	1.1 R	1.5 U	5110
Silver	0.39 R	0.52	0.2	0.46 R	0.47	5110
Thallium	0.73 U	0.79 U	0.72 U	0.83 U	1.2 U	67.5
Vanadium	17	13	21	23	26	1020
Zinc	29	24	19	49	89 🐩	100000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SS - Surface soil

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

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TABLE 3
SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

			<del>,                                     </del>				т	<u> </u>	
	GS-01-SS	GS-02-SS	GS-03-SS	GS-04-SS	GS-05-SS	GS-06-SS	GS-07-SS	GS-08-SS	PRG
ANALYTE	Background				 				
SVOC (ug/kg)									
1,1-Biphenyl	380 U	430 U	430 U	430 U	390 U	450 U	45 J	1400 U	23300000
2-Methylnaphthalene	380 U	430 U	430 U	430 U	390 U	450 U	120 J	1400 U	NL
Acenaphthene	380 U	430 U	430 U	430 U	390.U	450 U	170 J	1400 U	29200000
Acenaphthylene	380 U	430 U	430 U	430 U	120 J	450 U	320 J	480 J	NL .
Anthracene	380 U	430 U	430 U	430 U	470	450 U	970	930 J	100000000
Benzo(a)Anthracene	380 U .	430 U	430 U	430 U	77.0	450 U	2500	1500	2110
Benzo(b)Fluoranthene	380 U	430 U	420 U	430 U	1800 J ∴ ⊴	450 U	∴4600 J.	211000 J €	2110
Benzo(ghi)Perylene	380 U	430 U	430 U	430 U	380 J	450 U	690 ∰ ∜	33800	NL
Benzo(k)Fluoranthene	380 U	430 U	430 U	430 U	₽∂2100 Jar	450 U	<b>第 5200知識於</b>	章12000 J	21100
Benzo(a)Pyrene	380 U	430 U	430 U	430 U	810	450 U	2100 ₽ 7	3700	211
Carbazole	380 U	430 U	430 U	430 U	200 J	450 U	. ∌510 ⊱ i.J.	190 J	86200
Chrysene	380 U	430 U	430 U	430 U	<b>3-1500</b>	450 U	4900	2300	211000
Dibenzo(a,h)Anthracene	380 UJ	430 UJ	430 U	430 U	190 J	450 U	520	1700	211
Dibenzofuran	380 U	430 U	430 U	430 U	390 U	450 U	240 J	1400 U	1560000
Fluoranthene	380 U	430 U	430 U	430 U	2600	450 U	8200	2600	22000000
Fluorene	380 U	430 U	430 U	430 U	390 U	450 U	180· J	1400 U	26300000
Indeno (1,2,3-cd) Pyrene	380 UJ	430 UJ	430 U	430 U	610	450 U	1100	5000	2110
Naphthalene	380 U	430 U	430 U	430 U	390 U	450 U	87 J	1400 U	188000
Pentachlorophenol	970 UJ	1100· UJ	1100 U	1100 U	· 980 U	1100 U	980 U	3500 U	9000
Phenanthrene	380 U	430 U	430 U	430 U	320 J	450 U	<u>-</u> 4200	610 J	NL
Pyrene	380 U	430 U	430 U -	430 U	1500	450 U	6200	2500	29100000
Pesticides (ug/kg)				-					
4,4'-DDE (p,p'-DDE)	3.8 U	4.3 U	4.3 U	4.3 U	9.9 U	4.5 U	6.4 U	13 U	7020
4,4'-DDT (p,p'-DDT)	3.8 U	4.3 U	4:3 U	4.3 U	4 U	4.5 U	12 J	20/-41/J-13/8	7020
alpha-BHC	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U	359
Endrin	3.8 U	4.3 U	4.3 U	4.3 U	12 U	4.5 U	7.1 U	10 U	185000
Endrin Aldehyde	3.8 U	4.3 U	4.3 U	4.3 U	18 J	4.5 U	3.9 U	17 J	NL
Endrin Ketone	3.8 U	4.3 U	4.3 U	4.3 U	17 U	4.5 U	20 U	24 U	NL
Methoxychlor	20 U	22 U	22 U	22 U	48 U	23 U	47 U	74 NJ	3080000

### TABLE 3 SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-02-SS	GS-03-SS	GS-04-SS	GS-05-SS	GS-06-SS	GS-07-SS	GS-08-SS	PRG
ANALYTE	Background			e					<u> </u>
VOC (ug/kg)									
Acetone	100 J	180 J	25 J	28 J	91 J	82 J	87 J	74 J	54300000
Benzene	11 U	12 U	13 U	12 U	10 U	13 U	10 U	11 U	1410
Methyl Acetate	11 U	12 U	13 U	12 U	10 U	13 U	10 U	11 U	91500000
Methyl Ethyl Ketone	11 U	* 21	13 UJ	12 UJ	10 UJ	13 UJ	10 UJ	11 UJ	113000000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SS - Surface soil

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

N -

Presumptive evidence that analyte is present; reported as a tentative identification with an

NJ - estimated value.

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

SVOC - Semivolatile organic compounds

U - Analyte was not detected at or above the detection limit

μg/kg - Micrograms per kilogram

VOC - Volatile organic compounds

Samples were collected from 6 to 12 inches below ground surface; therefore, are considered surface soil for

\* - this investigation

TABLE 3
SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

·		<del></del>	<u> </u>		<u> </u>		<del></del>	1	·
	GS-01-SS	GS-09-SS	GS-10-SB*	GS-10-SS	GS-11-SS	GS-12-SS	GS-13-SS	GS-14-SB*	PRG
ANALYTE	Background_								
SVOC (ug/kg)									
1,1-Biphenyl	380 U	450 U	800 U	380 U	350 U	400 U	1100 U	390 U	23300000
2-Methylnaphthalene	380 U	450 U	200 J	69 J	70 J	400 U	1100 U	390 U	NL
Acenaphthene	380 U	450 U	210 J	97 J	350 U	400 U	1100 U	390 U	29200000
Acenaphthylene	380 U	450 U	<b>2</b> 61.00⊋	1100	330 J	400 U	1000 J	330 J	NL
Anthracene	380 U	450 U	<b>⊭12000</b>	1700	830	400 U	1600	690	100000000
Benzo(a)Anthracene	380 U	450 U	<b>≩23000</b>	2800	1000	400 U	4300	1400****	2110
Benzo(b)Fluoranthene	380 U	450 U	<b>\$37000</b>	5200	2500	400 U	7300 J	2400	2110
Benzo(ghi)Perylene	380 U	450 U	綠5200	1300	460	400 U	2000	540	NL
Benzo(k)Fluoranthene	380 U	450 U	26000	2300	∴ 1 <del>/</del> 700 ≦	. 400 U	5400 J	2000	21100
Benzo(a)Pyrene	380 U	450 U	25000	3100	<b>经</b> 素800	400 U	3100	1400	211
Carbazole	380 U	450 U	1800	- 480 #£,55	210 J	400 U	540 J	210 J	86200
Chrysene	380 U	450 U	35000	4300 金銭	1500	400 U	6200	2100	211000
Dibenzo(a,h)Anthracene	380 UJ	450 U	3300%	750	310 J	400 U	#1100	390	211
Dibenzofuran	380 U	450 U	340 J	120 J	140 J	400 U	1100 U	390 U	1560000
Fluoranthene	380 U	450 U	28000	4700	1900 E	400 U	11000	2200	22000000
Fluorene	380 U	450 U	800 U	380 U	350 U	400 U	1100 U	390 U	26300000
Indeno (1,2,3-cd) Pyrene	380 UJ	450 U	20000	2300	940	400 U	3100	1100	2110
Naphthalene	· 380 U	450 U	390 J	110 J	120 J	400 U	1100 U	390 U	188000
Pentachlorophenol	970 UJ	1100 U	680 J	690 J	640 J	1000 U	2900 U	980 U	9000
Phenanthrene	380 U	450 U	2700	1400	540	400 U	1600	\$510 <b>\$</b> \$	NL ·
Pyrene	380 U	450 U	., 37000 🗼 🗼 🛔	5000	1700	400 U	6800	2400	29100000
Pesticides (ug/kg)					··· = ···				
4,4'-DDE (p,p'-DDE)	3.8 U	4.5 U	4 U	3.8 U	3.5 U	4 U	30 U	3.9 U	7020
4,4'-DDT (p,p'-DDT)	3.8 U	4.5 U	110 N	43 J	# 0.30 J	4 U	△ 38 J	16·J壽25	7020
alpha-BHC	2 U	2.3 U	2 U	1.9 U	1.8 U	2 U	2 U	2 U	359
Endrin	3.8 U	4.5 U	37 U	15 U	3.5 U	4 U	42 J	3.9 U	185000
Endrin Aldehyde	3.8 U	4.5 U	15 U	3.8 U	3.5 U	4 U	49 J	3.9 U	NL
Endrin Ketone	3.8 U	4.5 U	180 N	3.8 U	3.5 U	4 U	150	3.9 U	NL
Methoxychlor	20 U	23 U	360 U	81 NJ	54 U	20 U	20 U	43 U	3080000

### TABLE 3 SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-09-SS	GS-10-SB*	GS-10-SS	GS-11-SS	GS-12-SS	GS-13-SS	GS-14-SB*	PRG
ANALYTE	Background		· 						
VOC (ug/kg)							• '		
Acetone	100 J	13 UJ	78 J	84 J	220 J	210 UJ	200 J	55 J	54300000
Benzene .	11 U	13 U	2 J	11 U	11 U	10 U	15 U	10 U	1410
Methyl Acetate	11 U	13 U	13 U	11 U	11 U	10 U	15 U	10 U	91500000
Methyl Ethyl Ketone	11 U	13 UJ	13 U	.11 U	21 J	14 J	15 J	10 U	113000000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SS - Surface soil

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

N -

Presumptive evidence that analyte is present; reported as a tentative identification with an

NJ - estimated value.

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PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

SVOC - Semivolatile organic compounds

U - Analyte was not detected at or above the detection limit

µg/kg - Micrograms per kilogram

VOC - Volatile organic compounds

Samples were collected from 6 to 12 inches below ground surface; therefore, are considered surface soil for

\* - this investigation

TABLE 3
SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

		<del> </del>	<del>r</del>	<del>1 :</del>	<del></del>		T	T	<del></del>
·	GS-01-SS	GS-14-SS	GS-15-SS	GS-16-SS	GS-17-SS	GS-18-SS	GS-19-SS	GS-20-SS	PRG
ANALYTE	Background	<u> </u>							
SVOC (ug/kg)		-						-	
1,1-Biphenyl	380 U	380 U	390 U	420 U	420 U	430 U	. 400 U	420 U	23300000
2-Methylnaphthalene	380 U	380 U	390 U	420 U	420 U	430 U	400 U	420 U	NL
Acenaphthene	380 U	380 U	390 U	420 U	420 U	430 U	400 U	420 U	29200000
Acenaphthylene	380 U	380 U	250 J	87 J	270 J	430 U	400 U	420 U	NL
Anthracene	380 U	1100	370 J	420 U	380 J	430 U	400 U	420 U	100000000
Benzo(a)Anthracene	380 U	2300	690c	260 J	1000	430 U	400 U	58 J	2110
Benzo(b)Fluoranthene	380 U	4200	∜ 1500‴₩¥X	700 J	2600 J ₹	430 U	400 U	160 J	2110
Benzo(ghi)Perylene	380 U	<b>7</b> 820	360 J	92 J	290 J	430 U	400 U	420 U	NL
Benzo(k)Fluoranthene	380 U	2400	1100	# 730 J	2700 J	430 U	400 U	170 J	21100
Benzo(a)Pyrene	380 U	<b>2300</b>	650	230 J	1000	430 U	400 U	46 J	211
Carbazole	380 U	280 J	250 J	62 J	130 J	430 U	400 U	420 U	86200
Chrysene	380 U	3100	1700	390 J	1600	430 U	400 U	98 J	211000
Dibenzo(a,h)Anthracene	380 UJ	580.	200 J	95 J	330 J	430 U	400 U	420 U	211
Dibenzofuran	380 U	380 U	390 U	420 U	420 U	430 U	400 U	420 U	1560000
Fluoranthene	380 U	3100	3100%	760	2500	430 U	400 U	73 J	22000000
Fluorene	380 U	380 U	390 U	420 U	420 U	430 U	400 U	420 U	26300000
Indeno (1,2,3-cd) Pyrene	380 UJ	1700	740 J	270 J	890	430 U	400 U	.63 J	2110
Naphthalene	380 U	380 U	390 U	420 U	420 U	430 U	400 U	420 U	188000
Pentachlorophenol	970 UJ	970 U .	990 UJ	1100 U	1100 U	1100 U	1000 U	1100 U	9000
Phenanthrene	380 U	320 J	<b>第1100 浅菜</b> 餐	310 J	250 J	430 U	400 U	420 U	NL
Pyrene	380 U	3300	。2200 清章	540	1700	430 U	400 U	. 100 J	29100000
Pesticides (ug/kg)					· · · · · · · · · · · · · · · · · · ·				
4,4'-DDE (p,p'-DDE)	3.8 U	3.8 U	3.9 U	2.5 J	4.2 U	4.3 U	4 U	4.1 NJ	7020
4,4'-DDT (p,p'-DDT)	3.8 U	10 NJ	30 NJ	4.2 U	4.2 U	4.3 U	4 U	4.2 U	7020
alpha-BHC	2 U	2 U	2 U	2.2 U	2.2 U	2.2 U	2 U	2.2 U	359
Endrin	3.8 U	3.8 U	11 U	4.2 U	6.4 J	4.3 U	4 U	4.2 U	185000
Endrin Aldehyde	3.8 U	3.8 U	27 NJ	4.2 U	4.2 U	4.3 U	4 U	4.2 U	NL
Endrin Ketone	3.8 U	3.8 U		14-3		4.3 U	4 U	4.2 U	NL
Methoxychlor	20 U	34 NJ	61 U	22 U	22 U	22 U	20 U	22 U	3080000

### TABLE 3 SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-14-SS	GS-15-SS	GS-16-SS	GS-17-SS	GS-18-SS	GS-19-SS	GS-20-SS	PRG
ANALYTE	Background	·							<u> </u>
VOC (ug/kg)									. •
Acetone	100 J	75 J	100 J	56 J	260 J	130 J	180 J	83 J	54300000
Benzene	11 U	11 U	10 U	12 U	13 U	13 U	11 U	12 U	1410
Methyl Acetate	11 U	11 U	3 J	12 U	13 U	13 U	3 J	12 U	91500000
Methyl Ethyl Ketone	11 U	11 U	12	12 U	28 J	14 J	17 J	12 U	113000000.

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SS - Surface soil

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

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Presumptive evidence that analyte is present; reported as a tentative identification with an

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µg/kg - Micrograms per kilogram

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Samples were collected from 6 to 12 inches below ground surface; therefore, are considered surface soil for

\* - this investigation

TABLE 3
SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

		<del> </del>	<u> </u>		1	
	GS-01-SS	GS-21-SS	GS-22-SS_	GS-23-SS	GS-24-SS	PRG
ANALYTE	Background			_		
SVOC (µg/kg)				_		•
1,1-Biphenyl	380 U	420 U	380 U	450 U	600 U	23300000
2-Methylnaphthalene	380 U	420 U	380 U	450 U	600 U	NL
Acenaphthene	380 U	420 U	380 U	450 U	600 U	29200000
Acenaphthylene	380 U	420 U	380 U	66 J	150 J	NL
Anthracene	380 U	420 U	380 U	430 U	330 J	100000000
Benzo(a)Anthracene	380 U	420 U	380 U		990	2110
Benzo(b)Fluoranthene	380 U	420 U	380 U	. NA	2700 J	2110
Benzo(ghi)Perylene	380 U	420 U	380 U	180 J	330 J	NL
Benzo(k)Fluoranthene	380 U	420 U	380 U	520	2600 J 🛂	21100
Benzo(a)Pyrene	380 U	420 U	380 U	380 J	870	211
Carbazole	380 U	420 U	380 U	450 U	600 U	86200
Chrysene	380 U	420 U	380 U	510	1100	211000
Dibenzo(a,h)Anthracene	380 UJ	420 U	380 U	120 J	260 J	211
Dibenzofuran	380 U	420 U	380 U	450 U	600 U	1560000
Fluoranthene	380 U	420 U	380 U	520	1300	22000000
Fluorene	380 U	420 U	380 U	450 U	600 U	26300000
Indeno (1,2,3-cd) Pyrene	380 UJ	420 U	.380 U	340 J	550 J	2110
Naphthalene	380 U	420 U	380 U	450 U	600 U	188000
Pentachlorophenol	970 UJ	1100 U	950 U	1100 UJ	1500 U	9000
Phenanthrene	380 U	420 U	380 U	59 J	600 U	NL
Pyrene	380 U	420 U	380 U	450	1600	29100000
Pesticides (ug/kg)					· ,	
4,4'-DDE (p,p'-DDE)	3.8 U	4.2 U	3.8 U	4.5 U	6 U	7020
4,4'-DDT (p,p'-DDT)	3.8 U	4.2 U	3.8 U	4.5 U	6 U	7020
alpha-BHC	2 U	2.2 U	2 U	1.7 NJ	3.1 U	359
Endrin	3.8 U	4.2 U	3.8 U	4.5 U	6 U	185000
Endrin Aldehyde	3.8 U	4.2 U	3.8 U	4.5 U	6 U	NL
Endrin Ketone	3.8 U	4.2 U	3.8 U	6.9 J	6 U	NL
Methoxychlor	20 U	22 U	20 U	23 U	31 U	3080000

## TABLE 3 SUMMARY OF 2003 SURFACE SOIL ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SS	GS-21-SS	GS-22-SS	GS-23-SS	GS-24-SS	PRG
ANALYTE	Background					
VOC (ug/kg)						
Acetone	100 J	270 J	110 J	130 J	220 J	54300000
Benzene	11 U	12 U	11 U	13 U	25 U	1410
Methyl Acetate	11 U	12 U	11 U	13 U	25 U	91500000
Methyl Ethyl Ketone	11 U	🐧 31 J 👔 🤻	11 UJ	14	25 UJ	113000000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SS - Surface soil

SB - Subsurface soil

GS - Gulf States Creosoting Company

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Presumptive evidence that analyte is present; reported as a tentative identification with an estimated

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µg/kg - Micrograms per kilogram

VOC - Volatile organic compounds

Samples were collected from 6 to 12 inches below ground surface; therefore, are considered surface soil

\* - for this investigation

TABLE 4
SUMMARY OF 2003 SUBSURFACE SOIL INORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

	GS-01-SB	GS-02-SB	GS-03-SB	GS-04-SB	GS-05-SB	GS-06-SB	GS-07-SB	GS-08-SB	PRG
ANALYTE	Background						· · · ·		TRO
Metals, Total (mg/kg)					•				-
Aluminum	7300 J	7800 J	8600 J	6600 J	5300 J	7700 J	9400 J	8100 J	100000
Arsenic	6.7	3.4 J	9.4	1.9	0.91 U	2.7	2.7	3.5	1.59
Barium	62	_ 34	63	57	59	48	50	. 45	66600
Beryllium	0.31	0.36	0.64	0.41	0.27	0.24	0.35	0.59	1940
Chromium	12	9.5	14	6.6	4.9	7.5	8	9.2	448
Cobalt	3.3	2.6	4.3	1.4	0.78 R	1.2	1.6	9.4	1920
Copper	6.4	3.8 UJ	6 U	4 U	2.1 U	3.4 U	5.7 U	3.9 U	40900
Iron	13000	13000	22000 J	7500 J	2300 J	5800 J	8200 J	12000 J	100000
Lead	8.2 J	6.6	15	9.2	6.7	11	11	5.7	800
Magnesium	580 J	480 J	770	370	220	580	550	370	NL
Manganese	180	200	200	60	16	18	19	-1000	19500
Selenium	0.7	1.4	1.6	0.66 U	0.62 U	0.96 U	0.65 U	0.58 U	5110
Silver	0.38	· 0.38 R	0.49 R	0.19 U	0.18 U	0.28 R	0.19 U	0.32	5110
Vanadium	22	20	33	14	6.7	17	18	22	1020
Zinc	18	20	23	15	7.3	12	. 15 .	25	100000

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

TABLE 4
SUMMARY OF 2003 SUBSURFACE SOIL INORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

	GS-01-SB	GS-09-SB	GS-11-SB	GS-12-SB	GS-13-SB	GS-15-SB	GS-16-SB	GS-17-SB	PRG .
ANALYTE	Background								
Metals, Total (mg/kg)	•								
Aluminum	7300 J	5800 J	9600 J	3700 J	8400 J	11000 J	12000 J	6700 J	100000
Arsenic	6.7	0.88 U	6.6	4.6	4.7	8.1	7.3	1.5	1.59
Barium	62	29	41	19	23	45	49	61	66600
Beryllium	0.31	0.35	0.49	0.3	0.34	0.59	0.49	0.39	1940
Chromium	12	5.3	17	8.5	11	21	. 17	7	448
Cobalt	3.3	1 R	3.2 R	1.4	1.7	2.9	2.1	1.5	1920
Copper	6.4	1.7 UJ	5.5 UJ	1.8 U	5.5 UJ	7 J	7.7	2.4 U	40900
Iron	13000	3800	20000	13000 J	14000	24000	23000	5900	100000
Lead	8.2 J	8.4	6.7	6.7	5.1	9.2	8.7 J	9.1 J	800
Magnesium	580 J	290 J	680 J	170	470 J	640 J	650 J	310 J	NL
Manganese	. 180	9.3	120	55	82	410	52	57	19500
Selenium	0.7	0.6 U	1.1 R	1.4	0.57 U	2.3 J	1.7 J	0.6 U	5110
Silver	0.38	0.18 U	1.1 R	0.47	0.39	0.65	0.45	0.32 R	5110
Vanadium	22	8.4	33	26	22	36	33	13	1020
Zinc	18	12	44	8.4	20	36	25	12 .	100000

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

TABLE 4
SUMMARY OF 2003 SUBSURFACE SOIL INORGANIC RESULTS
GULF STATES CREOSOTING COMPANY

	GS-01-SB	GS-18-SB	GS-19-SB	GS-20-SB	GS-21-SB	PRG
ANALYTE	Background					
Metals, Total (mg/kg)	<del>-</del> "					
Aluminum	7300 J	7700 J	12000 J	4700 J	8600 J	100000
Arsenic	6.7	4.1	8.3	2.3	4.5	1.59
Barium	62	26	43	24	28	66600
Beryllium	0.31	0.31	0.46	0.3	0.3	1940
Chromium	12	8.7	12	4.7	14	448
Cobalt	3.3	3.4	9.6	2.2	1.9	1920
Copper	6.4	3.9 U	6.3 U	2.3 U	3.6	40900
Iron	13000	12000	20000	6900	13000	100000
Lead	8.2 J	8 J	12 J	6 J	6.8 J	800
Magnesium	580.J	. 470 J	680 J	210 J	420 J	NL
Manganese	180	76	600	140	95	19500
Selenium	0.7	0.59 U	1.7 J	0.57 U	↑ 0.95 R	5110
Silver	0.38	0.32 R	0.53	0.25 R	0.4	5110
Vanadium	22	18	31	14 ·	21	1020
Zinc	18	16	30	11	16	100000

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

SB - Subsurface soil

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

R - Data is unusable

## TABLE 5 SUMMARY OF 2003 SUBSURFACE SOIL ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SB	GS-02-SB	GS-03-SB	GS-04-SB	GS-05-SB	GS-06-SB	GS-07-SB	GS-08-SB	PRG
ANALYTE	Background								
SVOC (ug/kg)						•			
Anthracene	400 U	410 U	460 U	430 U	410 U	450 U	450 U	74 J	100000000
Benzo(a)Anthracene	400 U	410 U	460 U	430 U	410 U	450 U	450 U	120 J	2110
Benzo(b)Fluoranthene	400 U	410 U	460 U	430 U	410 U	450 U	47 J	1000 J	2110
Benzo(ghi)Perylene	400 U	410 U	460 U	430 U	410 U	450 U	450 U	320 J	NL
Benzo(k)Fluoranthene	400 U	410 U	460 U	430 U	410 U	450 U	45 J	980 J 🤭 💰	21100
Benzo(a)Pyrene	400 U	410 U	460 U	430 Ü	410 U	450 U	450 U	240 J	211
bis(2-Ethylhexyl) Phthalate	400 U	410 U	460 UJ	430 UJ	450 J	450 U	450 U	400 U	123000
Chrysene	400 U	410 U	460 U	430 U	410 U	450 U	450 U	220 J	211000
Dibenzo(a,h)Anthracene	400 U	410 UJ	460 U	430 U	410 U	450 U	450 U	150 J	211
Fluoranthene	400 U	410 U	460 U	430 U	410 U	450 U	450 U	200 J	22000000
Indeno (1,2,3-cd) Pyrene	400 U	410 UJ	460 U	430 U	410 U	450 U	450 U	420	2110
Pyrene	400 U	410 U	460 U	430 U	410 U	450 U	450 U	230 J	29100000
Pesticides (ug/kg)									
4,4'-DDE (p,p'-DDE)	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U	7020
gamma-BHC (Lindane)	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U	1740
PCB-1260 (Aroclor 1260)	40 U	41 U	46 U	43 U	41 U	45 U	45 U	40 U	NL
VOC (ug/kg)									
Acetone	39 UJ	12 J	35 J	15 UJ	12 UJ	21 J	13 UJ	99 J 🕸	54300000
Methyl Ethyl Ketone	11 UJ	11 U	23 UJ	15 ŪJ	12 UJ	. 13 UJ	13 UJ	11 J	113000000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

a - Due to NJ qualification, compound was not used in HRS scoring

SB - Subsurface soil

GS - Gulf States Creosoting Company

HRS - Hazard Ranking System

J - Constituent was detected, the reported value is an estimate

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

SVOC - Semivolatile organic compounds

U - Analyte was not detected at or above the detection limit

μg/kg - Micrograms per kilogram

## TABLE 5 SUMMARY OF 2003 SUBSURFACE SOIL ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-SB	GS-09-SB	GS-11-SB	GS-12-SB	GS-13-SB	GS-15-SB	GS-16-SB	GS-17-SB	PRG
ANALYTE	Background								
SVOC (ug/kg)									
Anthracene	400 U	420 U	49 J	450 U	390 U	420 U	420 U	420 U	100000000
Benzo(a)Anthracene	400 U	420 U	400 U	450 U	56 J	420 U	420 U	420 U	. 2110
Benzo(b)Fluoranthene	400 U	420 U	400 U	450 U	160 J	420 U	420 U	41 J	2110
Benzo(ghi)Perylene	400 U	420 U	400 U	450 U	390 U	420 U	420 U	420 U	NL
Benzo(k)Fluoranthene	400 U	420 U	46 J	450 U	170 J	420 U	420 U	43 J	21100
Benzo(a)Pyrene	400 U	420 U	400 U	450 U	390 U	420 U	420 U	420 U	211
bis(2-Ethylhexyl) Phthalate	400 U	420 U	400 U	450 U	390 U	420 U	420 U	420 U	123000
Chrysene	400 U ·	420 U	400 U	450 U	91 J	420 U	420 U	420 U	211000
Dibenzo(a,h)Anthracene	400 U	420 U	400 U	450 U	390 U	420 UJ	420 U	420 U	· 211
Fluoranthene	400 U	420 U	400 U	450 U	130 J	420 U	420 U	420 U	22000000
Indeno (1,2,3-cd) Pyrene	400 U	420 U	400 U	450 U	51 J	420 UJ	420 U	420 U	2110
Pyrene	400 U	420 U	400 U	450 U	89 J	420 U	420 U	420 U	29100000
Pesticides (ug/kg)					*				
4,4'-DDE (p,p'-DDE)	4 U	4.2 U	4 U	4.5 U	6.9	4.2 U	4.2 U	. 4.2 U	7020
gamma-BHC (Lindane)	2.1 U	2.2 U	2 U	2.3 U	2 U	2.2 U	1.3 J	2.2 U	1740
PCB-1260 (Aroclor 1260)	40 U	42 U	40 U	45 U	39 U	42 U	<b>5.3</b> 70,	42 U	NL
VOC (ug/kg)									
Acetone	39 UJ .	11 UJ	24 J	11 UJ	21 UJ	67 J	11 UJ	13 U	54300000
Methyl Ethyl Ketone	11 UJ	11 UJ	11 UJ	11 UJ	10 UJ	11 U	11 UJ	13 U	113000000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

a - Due to NJ qualification, compound was not used in HRS scoring

SB - Subsurface soil

GS - Gulf States Creosoting Company

HRS - Hazard Ranking System

J - Constituent was detected, the reported value is an estimate

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

SVOC - Semivolatile organic compounds

U - Analyte was not detected at or above the detection limit

µg/kg - Micrograms per kilogram

## TABLE 5 SUMMARY OF 2003 SUBSURFACE SOIL ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

·	GS-01-SB	GS-18-SB	GS-19-SB	GS-20-SB	GS-21-SB	PRG
ANALYTE	Background					FIG
SVOC (ug/kg)						
Anthracene	400 U	440 U	420 U	400 U	410 U ·	100000000
Benzo(a)Anthracene	400 U	440 U	420 U	400 U	410 U	2110
Benzo(b)Fluoranthene	400 U	440 U	420 U	400 U	410 U	2110
Benzo(ghi)Perylene	400 U	440 U	420 U	120 J	410 U	NL
Benzo(k)Fluoranthene	400 U	440 U	420 U	400 U	410 U	21100
Benzo(a)Pyrene	400 U	440 U	420 U	400 U	410 U	211
bis(2-Ethylhexyl) Phthalate	400 U	440 UJ	420 UJ	400 U	410 U	123000°
Chrysene	400 U	440 U	420 U	400 U	410 U	211000
Dibenzo(a,h)Anthracene	400 U	440 U	420 U	400 U	410 U	211
Fluoranthene	400 U	440 U	420 U	400 U	410 U	22000000
Indeno (1,2,3-cd) Pyrene	400 U	440 U	420 U	400 U	410 U	2110
Pyrene	400 U	440 U	420 U	400 U	410 U	29100000
Pesticides (ug/kg)				-		
4,4'-DDE (p,p'-DDE)	4 U	4.4 U	4.2 U	4 U	4.1 Ü	7020
gamma-BHC (Lindane)	2.1 U	1.7 J	2.2 U	2.1 U	2.1 U	1740
PCB-1260 (Aroclor 1260)	40 U	42 J	39 J	40 U	41 U	NL
VOC (ug/kg)						
Acetone	39 UJ	14 UJ	26 UJ	11 U	11 UJ	54300000
Methyl Ethyl Ketone	11 UJ	11 UJ	11 UJ	11 U	11 UJ	113000000

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds PRG

a - Due to NJ qualification, compound was not used in HRS scoring

SB - Subsurface soil

GS - Gulf States Creosoting Company

HRS - Hazard Ranking System

J - Constituent was detected, the reported value is an estimate

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Industrial Soil

SVOC - Semivolatile organic compounds

U - Analyte was not detected at or above the detection limit

μg/kg - Micrograms per kilogram

## TABLE 6 GULF STATES CREOSOTING COMPANY SUMMARY OF 2003 GROUNDWATER INORGANIC RESULTS

	GS-01-GW	GS-02-GW	GS-03-GW	GS-04-GW	GS-05-GW	MCL	PRG
ANALYTE	Background						FRG
Metals, Total (ug/L)							
Aluminum	320 UJ	450°J	₹ 850 J	1700 J	370°J, 🕾 👙	. NL	36500
Barium	20	\$ \$120\cdots	29	26	32	2000	2550
Beryllium	0.11	0.14	0.11	0.1 U	0.9	4	73
Chromium	1.8 R	0.92	1.8	1.6	0.6 U	100	NL
Cobalt	1.2	1.2 R	0.9 U	0.9 U	10	NL	730
Copper	1.5 U	0.73 U	1.1 U	0.78 U	0.6 U	1300	1460
Iron	1500	1400	2100	1900	12000	NL	10900
Magnesium	1500	5500	940	930	8200	NL	NL
Manganese	23	160	<b>第 3110</b> 年(春)	. 30	320	NL	876
Nickel	1.5 U	3.5	1.5 U	1 14 3.31 A	⊋20 <b>≨</b>	NL	730
√anadium	0.6 U	0.82 R	2 0 V	第4-1.9 0	0.6 U	NL	36.5
Zinc	11 U	132	2 26 18 20 8	11 U	**************************************	NL	10900

#### Notes:

Shaded - Concentration is elevated (3 x background)

MCL - EPA Maximum Contaminant Limit

Bold - Value exceeds PRG

GW - Groundwater sample

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

ug/L - Micrograms per liter

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Tap Water

R - Data is unusable

## TABLE 7 SUMMARY OF 2003 GROUNDWATER ORGANIC RESULTS GULF STATES CREOSOTING COMPANY

	GS-01-GW	GS-02-GW	GS-03-GW	GS-04-GW	GS-05-GW	MCL	PRG
ANALYTE	Background						
VOC (ug/L)							
Methyl Ethyl Ketone	6.1 UJ	5 UJ	6.9 UJ	7:2 J	6.6 UJ	NL	6970

#### Notes:

Shaded - Concentration is elevated (3 x background)

MCL - EPA Maximum Contaminant Limit

GW - Groundwater sample

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

ug/L - Micrograms per liter

NL - No limit established

PRG - Region 9 Preliminary Remediation Goal for Tap Water

## TABLE 8 SUMMARY OF 2003 SEDIMENT INORGANIC RESULTS - MARSH SAMPLES GULF STATES CREOSOTING COMPANY

	GS-01-SD	GS-02-SD	GS-03-SD	GS-04-SD	GS-05-SD	EcoScreen
		00-02-00		GG-04-GD	- <del> </del>	Sediment
ANALYTE	Background				·	
Metals, Total (mg/kg)			:	÷		
Aluminum	8200 J	4700 J	16000 J	9100 J	4900 J	NL
Arsenic	5.7	1.8 R	7.9 R	5.5	2.9 R	7.24
Barium	89	77	ુ: 300 ખેત	120	71	NL
Beryllium	0.39	0.56	# #1.3. **.***	* -1.25°	0.51	NL
Chromium	13	7.8	24	15	8.4	52.3
Cobalt	4.4	4.7	12	11	5.5	NL
Iron	15000	6400	24000	13000	8400	NL
Lead	12 J	20 J	41 J	21	12	30.2
Magnesium	1000 J	410 J	1200 J	1000 J	660 J	NL
Manganese	340	500	720	280	320	NL
Selenium	1.9 R	1.1 U	4.2	2.1 U	- 1 U	NL
Vanadium	24	13	41	23	13	NL
Zinc	36	33	. 3130	84	39	124

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds screening level

SD - Sediment sample

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

NL - No limit established

R - Data is unusable

TABLE 10
SUMMARY OF 2003 SEDIMENT INORGANIC RESULTS - PEARL RIVER SAMPLES
GULF STATES CREOSOTING COMPANY

ANALYTE	GS-06-SD Control	GS-07-SD	EcoScreen Sediment
Metals, Total (	mg/kg)		
Aluminum	450 J	140 J	NL
Barium	6.8	3.2	NL
Beryllium	0.03 U	0.03 U	NL
Chromium	1.3 U	0.61 U	52.3
Cobalt	0.61 R	0.47	NL
Iron	950	340	NL
Lead	1.8	1.2 J	30.2
Magnesium	64 J	14 UJ	NL
Manganese	52	53	NL
Vanadium	1.6 R	0.57 R	NL
Zinc	3.6	2.9 U	124

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds screening level

SD - Sediment Sample

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

mg/kg - Milligrams per kilogram

NL - No limit established

R - Data is unusable

# TABLE 9 SUMMARY OF 2003 SEDIMENT ORGANIC RESULTS - MARSH SAMPLES GULF STATES CREOSOTING COMPANY

·	GS-01-SD	GS-02-SD	GS-03-SD	GS-04-SD	GS-05-SD	EcoScreen
ANALYTE	Background					Sediment
SVOC (ug/kg)	•					
Anthracene	440 U	410 U	270 J	900 U	430 U	330
Benzaldehyde	89 J	410 U	110 J	130 J	430 U	NL
Benzo(a)Anthracene	440 U	410 U	150 J	900 U	430 U	330
Benzo(b)Fluoranthene	440 U	410 U	540 J	900 U	430 U	655
Benzo(ghi)Perylene	92 J	410 U	940 U	900 U	430 U	655
Benzo(k)Fluoranthene	440 U	410 U	₹ 560 J	900 U	430 U	655
Benzo(a)Pyrene	440 U	410 U	150 J	900 U	430 U	330
Chrysene	440 U	410 U	250 J	900 U	430 U	330
Fluoranthene	440 U	410 U	340 J	900 U	430 U	330
Indeno (1,2,3-cd) Pyrene	440 U	410 U	180 J	900 U	430 UJ	655
Naphthalene	440 U	410 U	940 U	3 J	430 U	330
Phenanthrene	440 U	410 U	120 J	900 U	430 U	330
Pyrene	440 U	410 U	220 J	120 J	430 U	330
PAH (ug/kg)						
2-Methylnaphthalene	0.6 J	0.9 J	∜r + 33i	2 J	0.8 J	330
Acenaphthene	0.5 J	1 J	1,00%	5.	0.9 J	330
Acenaphthylene	3	. 2	44.	9	. 2 U	330
Anthracene	3	7	140	34	7	330
Benzo(a)Anthracene	1 J	<b>学生7 学</b> 漢	110	19	12	330
Benzo(a)Pyrene	2	4		# F3.17: 11:	16 <sup>1</sup>	330
Chrysene	1 J	10	200	36	. 26 F	330
Dibenzo(a,h)Anthracene	1 U	1 J	** ** 40	6	*4	330
Fluoranthene	2	21 %	190	50	34	330
Fluorene	0.5 J	2	. 82	6	· 1 J	330
Naphthalene	0.6 J	0.6 J	51	8 3 J	0.9 J	330
Phenanthrene	2 J	<b>12</b>	72	11	23 U	330
Pyrene	2	15,3	140	33	27	330

## TABLE 9 SUMMARY OF 2003 SEDIMENT ORGANIC RESULTS - MARSH SAMPLES GULF STATES CREOSOTING COMPANY

	GS-01-SD	GS-02-SD	GS-03-SD	GS-04-SD	GS-05-SD	EcoScreen
ANALYTE	Background				·	Sediment
Pesticides (ug/kg)						
4,4'-DDE (p,p'-DDE)	7.3 U	7.9	16 U	15 U	7.2 U	3.3
beta-BHC	3.8 U	3.5 U	12	7.7 U	3.7 U	NL
Endosulfan II (beta)	7.3 U	6.5 J	16 U	15 U	7.2 U	NL
Endosulfan Sulfate	7.3 U	10 N	14 NJ	15 U	7.2 U	NL
VOC (ug/kg)	•					
Acetone	94 J	130 J	≛ 350 ป	520 J	51 J	NL
Carbon Disulfide	5 J	34 U	ຸ 51∜ປ	70 U	29 U	NL
Methyl Ethyl Ketone	37 UJ	34 UJ	75 U	∴88%J	29 U	NL

#### Notes:

Shaded - Concentration is elevated (3 x background)

Bold - Value exceeds screening level

SD - Sediment Sample

GS - Gulf States Creosoting Company

J - Constituent was detected, the reported value is an estimate

N-

NJ - Presumptive evidence that analyte is present; tentative identification with an estimated value.

PAH - Polycyclic aromatic hydrocarbon

NL - No limit established

R - Data is unusable

SVOC - Semivolatile organic compounds

U - Analyte was not detected at or above the detection limit

μg/kg - Micrograms per kilogram

### TABLE 11 **SUMMARY OF 2003 SEDIMENT ORGANIC RESULTS - PEARL RIVER SAMPLES GULF STATES CREOSOTING COMPANY**

	GS-06-SD	GS-07-SD	EcoScreen
ANALYTE	. Control		Sediment
PAH (ug/kg)	•		
2-Methylnaphthalene	0.4 J	0.4 J	330
Dibenzo(a,h)Anthracene	1	0.8 U	330
Fluorene	0.3 J	0.8 U	330
Naphthalene	0.5 J	0.4 J	330
Phenanthrene	4	0.9 U	330
VOC (ug/kg)			
Acetone	22 J	49 J	NL

#### Notes:

GS - Gulf States Creosoting Company

SD - Sediment Sample

HRS - Hazard Ranking System

J - Constituent was detected, the reported value is an

estimate

PAH - Polycyclic aromatic hydrocarbon

NL - No limit established

U - Analyte was not detected at or above the detection

µg/kg - Micrograms per kilogram



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Transmitted Electronically

December 31, 2003

Ms. Donna Webster Remedial Project Manager U.S. Environmental Protection Agency, Region 4 61 Forsyth Street, SW, 11th Floor Atlanta, Georgia 30303

Subject:

Final Preliminary Assessment/Site Inspection Report, Revision 1

**Gulf States Creosoting** 

EPA ID No. MSN000407423 EPA Contract No. 68-W-00-123

Technical Direction Document (TDD) No. 4W-02-03-A-003

**Document Control Number (DCN) WSI-GSC-0011** 

Dear Ms. Webster:

Weston Solutions, Inc., Superfund Technical Assessment and Response Team - 2 (START-2) is submitting two copies of the final preliminary assessment/site inspection (PA/SI) report, Revision 1 for the Gulf States Creosoting facility in Flowood, Rankin County, Mississippi and an additional copy of references for the state.

Please contact me at (404) 527-7016 if you have any questions regarding this final PA/SI report.

Sincerely,

Weston Solutions, Inc.

Shanieka Pennamon

START-2 Project Manager

Enclosure

cc:

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Alexis K. Ullock, START-2 Site Assessment Coordinator (w/o enclosure)

START-2 File

### FINAL PRELIMINARY ASSESSMENT/SITE INSPECTION REPORT

### GULF STATES CREOSOTING COMPANY FLOWOOD, RANKIN COUNTY, MISSISSIPPI U.S. EPA ID NO. MSN000407423

#### **Revision 1**

#### Prepared for

## U.S. ENVIRONMENTAL PROTECTION AGENCY Region 4 Atlanta, Georgia 30303

Contract No. : 68-W-00-123

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**Gulf States Creosoting Company** 

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.1	General Site Location Map
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1.0 INTRODUCTION

The U.S. Environmental Protection Agency (EPA) tasked the Weston Solutions, Inc. Superfund Technical Assessment and Response Team - 2 (START-2) to prepare a preliminary assessment/site inspection (PA/SI) report for the Gulf States Creosoting Company (Gulf States) facility in Flowood, Rankin County, Mississippi, EPA ID No. MSN000407423. The PA/SI was conducted under Contract No. 68-W-00-123, Technical

Direction Document (TDD) No. 4W-02-03-A-003.

The primary objective of a PA/SI is to determine whether a site has the potential to be placed on the National Priorities List (NPL). The NPL identifies sites at which a release, or threatened release, of hazardous substances poses a serious enough risk to public health or the environment to warrant further investigation and possible remediation under the Comprehensive Environmental Response, Compensation, and Liability

Act of 1980 (CERCLA) and the Superfund Amendments and Reauthorization Act of 1986.

Information gathered during the PA/SI is used to generate a preliminary Hazard Ranking System (HRS) score. The HRS is the primary criterion EPA uses to determine whether a site should be placed on the NPL. PA/SIs are generally conducted at sites where environmental sampling and/or monitoring well installation are necessary to fulfill HRS documentation requirements.

Specifically, the objectives of the PA/SI are as follows:

- Obtain and review relevant file material
- Collect samples to attribute hazardous substances to site operations
- Collect samples to establish representative background levels
- Evaluate target populations for the groundwater migration, surface water migration, soil exposure, and air migration pathways
- Collect any other missing HRS data
- Document current site conditions
- Develop a site layout map

This report documents the results of the PA/SI conducted at the Gulf States facility during the week of April 21, 2003. All sampling was conducted by personnel from EPA's Science and Ecosystem Support Division (SESD) (Refs. 1; 2). Information reviewed for the PA/SI was gathered from the Mississippi Department of Environmental Quality (MDEQ) and from the EPA Region 4 CERCLA files.

This document was prepared by Weston Solutions, Inc., expressly for EPA. It shall not be disclosed, in whole or in part, without the express written permission of EPA.

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DCN; WSI-GSC-0011

#### 2.0 SITE BACKGROUND

This section describes the facility, its present and past operations (including waste disposal practices and regulatory history), previous investigations, and potential source areas located at the facility.

#### 2.1 SITE DESCRIPTION AND ENVIRONMENTAL SETTING

The former 141-acre Gulf States Creosoting Company is located at 1625 Flowood Drive (Mississippi Hwy 468), Flowood, Rankin County, Mississippi (Refs. 3; 4, p. v; 5, pp. 1-4). The geographic coordinates of the facility are 32°18'43.8" north latitude and 90°58'38.3" west longitude (Ref 6). The facility is currently operating as a horse farm and as ConSteel Co, Inc.(ConSteel). Both current operations appear to be located on portions of the property where creosoting operations were conducted (Ref. 11). Currently, the southernmost building on the property is owned by ConSteel. The northernmost structures (i.e. metal barn and shed) are used to house tractors and helicopters for a local news station and are part of the horse farm. The horse farm has been operating on the property for approximately 7 years. The facility is bound by railroad tracks to the north and east, an adjacent business to the south, and marshland/tributary of the Pearl River to the west. The facility is bound by natural barriers and is not fenced (Refs. 3; 5). The facility lies within an area comprised of mixed industrial, commercial, and residential uses (Ref. 3).

The climate of Rankin County is characterized by long, hot summers and mild winters. Moist tropical air from the Gulf of Mexico has a moderating influence on maximum temperatures in summer (Ref. 7, p. 1). Normal annual total precipitation for Rankin County is approximately 55 inches, and the mean annual lake evaporation is 44 inches, yielding a net annual precipitation of 10 inches (Ref. 8). The 2-year, 24-hour rainfall is 4.5 inches (Ref. 9).

#### 2.2 SITE OPERATIONS AND REGULATORY HISTORY

Gulf States owned the property as early as 1929 (Ref. 10, p. 11). Gulf States operated as a wood treating facility until the mid 1950s (Ref. 5, pp. 2, 3). In July 1958, American Creosoting Corporation obtained portions of the property (Ref. 10, pp. 7-11). In June 1959, W.G. Avery Body Company obtained portions of the property and operated a body shop on the property (Ref. 4, Deeds, p. 6). There are several tracts of

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land on the Gulf States property. The horse farm is owned and operated by Mr. Jim Webb; Webb owns the

portion of the property just north of the main entrance to the property. ConSteel owns a portion of the

property due south of the main entrance. ConSteel appears to be operating on the portion of the property that

it owns. The onsite levee is owned by the Levee Board - which is a joint venture between the state and

Rankin County. Mr. John McGowan owns the marsh area due west of the levee. McGowan has plans of

developing the marsh areas as residential subdivisions (Ref. 30).

During Gulf States operations, railroad cross ties were treated at the facility with coal-tar creosote and

transported on and off-site by means of railroad box cars (Refs. 5, p. 2; 11). Creosote is typically applied

to the wood by commercial pressure treatment or by home and farm dipping (Ref. 12, p. 69). Creosote is a

wood preservative used to treat railroad ties, telephone poles, marine pilings, and fence posts (Ref. 12, pp.

1, 2).

A Phase I investigation of the property in July 1993 included a review of the MDEQ Underground Storage

Tank (UST) division records. No past UST usage on the property was found during the review (Ref. 4, p.

xii). No other environmental permits have been identified for the facility.

2.3 PREVIOUS RELEASES AND INVESTIGATIONS

In June 1993, BCM Engineers, Inc. (BCM) conducted a Phase I Environmental Assessment (Phase I) of the

former Gulf States property on behalf of Trustmark National Bank. The objective of the Phase I was to

identify adverse environmental conditions, suspect activities, and potential hazardous wastes or materials

on or in the vicinity of the subject property. The Phase I included the collection of soil samples from soil

borings ranging in depth from 0 to 8 feet below ground surface (bgs). The seven soil borings were

concentrated in the suspected location of the former creosote operations. Of the seven borings, five samples

were collected and analyzed for semi-volatile organic compounds (SVOC). Constituents detected in the soil

samples were napthalene, 2-methylnapthalene, dibenzofuran, fluorene, phenanthrene, and fluoranthene. The

concentrations ranged from below the detection limit to 604 milligrams per kilogram (mg/kg) (Ref. 4).

In August 1993, BCM conducted a Phase II Environmental Assessment of the former Gulf States property.

The objective of the Phase II was to delineate the extent of the creosote contaminated soil identified during

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the Phase I investigation. During the Phase II, several soil borings were advanced and five soil samples were collected for SVOC laboratory analysis. The soil borings ranged in depth from 0 to 8 feet bgs. Sample results indicated the presence of 2-methylnapthalene, naphthalene, dibenzofuran, phenanthrene, fluoranthene, pyrene, and chrysene. Constituents ranged from below the detection limit to 1,057 mg/kg. The data were compared to health-based criteria for exposure via groundwater ingestion. Based on such a comparison, the recommendation for no further action was presented because the soil contaminants were below the target cleanup levels (Ref. 13).

#### 2.4 POTENTIAL SOURCE AREAS

The source considered for the purpose of this PA/SI is contaminated soil.

#### 3.0 PA/SI ACTIVITIES

This section outlines field observations and sampling procedures at the sampling locations. Individual subsections address the sampling investigation and rationale for specific PA/SI activities. The PA/SI was conducted in accordance with the EPA Quality Assurance Project Plan (QAPP) for the Gulf States facility (Ref. 2). The QAPP was prepared by SESD and includes the sample recommendations submitted by START-2 (Refs. 2; 14). The only deviation from the sampling plan involved the addition of a control sample along Prairie Branch in order to attribute contamination to another source upstream of the Gulf States property (Ref. 5, p.11).

#### 3.1 SAMPLE COLLECTION METHODOLOGY AND PROCEDURES

SESD personnel collected 24 surface soil samples, 20 subsurface soil samples, four groundwater samples, and five sediment samples on or near the Gulf States property during the week of April 21, 2003 (Ref. 5). Sample locations are depicted on Figure 1 in Reference 1, Appendix A and are summarized in Tables 1 through 4 of this report.

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SESD personnel collected the surface soil samples from 0 to 6 inches below ground surface (bgs), and subsurface soil samples were collected from 2 to 3 feet bgs; two of the subsurface soil samples were collected between 6 and 12 inches. SESD personnel followed sample collection procedures outlined in the EPA Region 4 SESD, *Environmental Investigation Standard Operating Procedures and Quality Assurance Manual* (EISOPQAM) and the *Analytical Support Branch Operations and Quality Control Manual* (November 2001).

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TABLE 1
SURFACE SOIL SAMPLING LOCATIONS AND RATIONALE

Sample Number	Location	Rationale
GS-01-SS	Background; from Jackson Prep School (3100 Lakeland Dr.) located northeast of the property	Background soil sample for comparison to on-site sample results
GS-02-SS	North-central portion of the property	Determine presence or absence of hazardous substances
GS-03-SS	Depression located on the west-central portion of the property, west of the levee	Determine presence or absence of hazardous substances
GS-04-SS	Southwestern portion of the property	Determine presence or absence of hazardous substances
GS-05-SS	Southeastern portion of the property	Determine presence or absence of hazardous substances
GS-06-SS	Southwestern portion of the property	Determine presence or absence of hazardous substances
GS-07-SS	Southeastern portion of the property	Determine presence or absence of hazardous substances
GS-08-SS	Southeastern portion of the property	Determine presence or absence of hazardous substances
GS-09-SS	Southwestern portion of the property	Determine presence or absence of hazardous substances
GS-10-SS	Southeastern portion of the property near the existing ConSteel building	Determine presence or absence of hazardous substances
GS-11-SS	Central portion of the property from the barren area west of the ConSteel building	Determine presence or absence of hazardous substances
GS-12-SS	Western portion of the property, west of the levee	Determine presence or absence of hazardous substances
GS-13-SS	Central portion of the property south of the barn	Determine presence or absence of hazardous substances

### TABLE 1 (Continued)

### SURFACE SOIL SAMPLING LOCATIONS AND RATIONALE

Sample Number	Location	Rationale
GS-14-SS	Central portion of the property west of the barn	Determine presence or absence of hazardous substances
GS-15-SS	East-central portion of the property	Determine presence or absence of hazardous substances
GS-16-SS	Northeastern portion of the property	Determine presence or absence of hazardous substances
GS-17-SS	Northeastern portion of the property	Determine presence or absence of hazardous substances
GS-18-SS	Northeastern portion of the property	Determine presence or absence of hazardous substances
GS-19-SS	North-central portion of the property	Determine presence or absence of hazardous substances
GS-20-SS	North-central portion of the property	Determine presence or absence of hazardous substances
GS-20-SD	North-central portion of the property; duplicate of GS-20-SS	Determine presence or absence of hazardous substances
GS-21-SS	Northwestern portion of the property	Determine presence or absence of hazardous substances
GS-22-SS	Northern portion of the on-site drainage ditch	Determine presence or absence of hazardous substances
GS-23-SS	Central portion of the on-site drainage ditch	Determine presence or absence of hazardous substances
GS-24-SS	Southern portion of drainage ditch near fenced area	Determine presence or absence of hazardous substances

Notes

GS Gulf States Creosoting Company

SD - Duplicate surface soil sample

SS Surface soil sample

Surface soil samples were collected from 0 to 6 inches below ground surface (bgs)

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TABLE 2
SUBSURFACE SOIL SAMPLING LOCATIONS AND RATIONALE

Sample Number	Location	Rationale
GS-01-SB	Background; from Jackson Prep School (3100 Lakeland Dr.) located northeast of the property	Background soil sample for comparison to on-site sample results
GS-02-SB	North-central portion of the property	Determine presence or absence of hazardous substances
GS-03-SB	Depression located on the west-central portion of the property, west of the levee	Determine presence or absence of hazardous substances
GS-04-SB	Southwestern portion of the property	Determine presence or absence of hazardous substances
GS-05-SB	Southeastern portion of the property	Determine presence or absence of hazardous substances
GS-06-SB	Southwestern portion of the property	Determine presence or absence of hazardous substances
GS-07-SB	Southeastern portion of the property	Determine presence or absence of hazardous substances
GS-08-SB	Southeastern portion of the property	Determine presence or absence of hazardous substances
GS-09-SB	Southwestern portion of the property	Determine presence or absence of hazardous substances
GS-10-SB*	Southeastern portion of the property near the existing ConSteel building	Determine presence or absence of hazardous substances
GS-11-SB	Central portion of the property from the barren area west of the ConSteel building	Determine presence or absence of hazardous substances
GS-12-SB	Western portion of the property, west of the levee	Determine presence or absence of hazardous sûbstances
GS-13-SB*	Central portion of the property south of the barn	Determine presence or absence of hazardous substances

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## TABLE 2 (Continued)

### SUBSURFACE SOIL SAMPLING LOCATIONS AND RATIONALE

Sample Number	Location	Rationale
GS-14-SB	Central portion of the property west of the barn	Determine presence or absence of hazardous substances
GS-15-SB	East-central portion of the property	Determine presence or absence of hazardous substances
GS-16-SB	Northeastern portion of the property	Determine presence or absence of hazardous substances
GS-17-SB	Northeastern portion of the property	Determine presence or absence of hazardous substances
GS-18-SB	Northeastern portion of the property	Determine presence or absence of hazardous substances
GS-19-SB	North-central portion of the property	Determine presence or absence of hazardous substances
GS-20-SB	North-central portion of the property	Determine presence or absence of hazardous substances
GS-21-SB	Northwestern portion of the property	Determine presence or absence of hazardous substances

Notes:

GS Gulf States Creosoting Company

SB Subsurface soil sample

\* Sample collected from 6 inches to 1 foot bgs.

Subsurface soil samples were collected from 2 to 3 feet bgs, except as noted.

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TABLE 3

GROUNDWATER SAMPLING LOCATIONS AND RATIONALE

Sample Number	Location	Rationale
GS-01-GW	Background; from Jackson Prep School (3100 Lakeland Dr.) located northeast of the property	Background groundwater sample for comparison to on-site sample results
GS-02-GW	North-central portion of the property	Determine presence or absence of hazardous substances
GS-03-GW	Depression located on the west- central portion of the property, west of the levee	Determine presence or absence of hazardous substances
GS-04-GW	Southwestern portion of the property	Determine presence or absence of hazardous substances
GS-05-GW	Southeastern portion of the property	Determine presence or absence of hazardous substances

Notes:

GS

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GW

Groundwater sample

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TABLE 4
SEDIMENT SAMPLING LOCATIONS AND RATIONALE

Sample Number	Location	Rationale
GS-01-SD	Background; from tributary to Pearl River accessed from Jackson Prep School (3100 Lakeland Dr.) located northeast of the property	Background sediment sample for comparison to downstream marsh sample results
GS-02-SD	Northern portion of tributary to the Pearl River; area north of gated dirt road leading to the tributaries	Determine presence or absence of hazardous substances in the marsh
GS-03-SD	Southern portion of tributary to the Pearl River; area south of gated dirt road leading to the tributaries	Determine presence or absence of hazardous substances in the marsh
GS-04-SD	Half-way to the Pearl River along the southern portion of the tributary to the Pearl River	Determine presence or absence of hazardous substances in the marsh
GS-05-SD	Confluence of the southern tributary to the Pearl River and its tributary	Determine presence or absence of hazardous substances in the marsh
GS-06-SD	Control from Pearl River; approximately 1/8 mile upstream of confluence of Pearl River and Prairie Branch	Control sample for comparison to on-site sample results
GS-07-SD	Pearl River, downstream of the confluence of the tributary adjacent to Gulf States Creosoting and the Pearl River	Determine presence or absence of hazardous substances
GS-08-SD	Prairie Branch	Control sample for comparison to on-site sample results

Notes:

GS Gulf States Creosoting Company

SD Sediment sample

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3.2 ANALYTICAL SUPPORT AND METHODOLOGY

All samples collected during the PA/SI were analyzed through the EPA Contract Laboratory Program (CLP).

The laboratories analyzed for EPA Target Compound List (TCL) volatile organic compounds (VOC),

extractable semivolatile organic compounds (SVOC), pesticides, and polychlorinated biphenyls (PCB).

Certain samples were analyzed for specific SVOC concentrations at lower concentrations than the standard

levels of analysis under the CLP. The samples were also analyzed for Target Analyte List (TAL) inorganic

substances (metals) and cyanide. EPA Region 4 SESD reviewed all data for compliance with the terms of

the CLP.

3.3 ANALYTICAL DATA QUALITY AND DATA QUALIFIERS

All analytical data were subject to a quality assurance review as described in the EPA SESD laboratory data

evaluation guidelines. The text and analytical data tables presented in this report show some concentrations

of organic and inorganic parameters as qualified with a "J," indicating that the qualitative analysis was

acceptable; however, the quantitative value has been estimated. Other compounds may have been qualified

with an "N," indicating that they were detected based on the presumptive evidence of their presence. This

means that the compound was only tentatively identified, and its detection cannot be considered a positive

indication of its presence. Some sample results are reported with a "U" qualifier, meaning that the material

was analyzed for but not detected. The reported number is the laboratory-derived sample quantitation limit

(SQL) for the constituent in that sample. At times, miscellaneous organic compounds that do not appear on

the TCL are reported with the data set. These constituents are qualified as "JN," indicating that they are

tentatively identified at estimated quantities. Because these constituents are not routinely analyzed for or

reported, background levels of SQLs are not generally available for comparison. Some compounds are

qualified with an "R" which indicates the QC evaluation has determined the concentration of the compound

is unusable. Compounds qualified with a "C" have been confirmed by gas chromatograph or mass

spectrometry. The complete analytical data sheets are presented in Reference 1.

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4.0 SOURCE SAMPLING

This section discusses the source area evaluated at the facility and the sampling locations and analytical

results of samples collected from the Gulf States property. The source area at the Gulf States property

evaluated in this PA/SI is contaminated soil...

Surface soil and subsurface soil sampling locations are depicted on the sample location map in Reference

1, Appendix A and described in Tables 1 and 2. Surface soil inorganic and organic analytical results are

summarized in Tables 5 and 6, respectively, and subsurface soil inorganic and organic analytical sampling

results are summarized in Tables 7 and 8, respectively. Tables 5 through 8 are presented following Section

6.0. Elevated concentrations of constituents are shaded in the tables. The concentration of a constituent is

considered to be elevated if the concentration is greater than or equal to three times the concentration

detected in the background or control sample. In the case where a constituent is undetected in the

background or control sample, any concentration equal to or greater than the SQL is considered to be

elevated. The summary analytical data tables are presented as Appendix A.

The following discussion of hazardous constituents detected at elevated levels in soil samples collected at

the facility includes only those hazardous constituents that are associated with site operations and those

hazardous constituents that may pose a threat to human health or the environment.

4.1 SOURCE SAMPLING LOCATIONS AND ANALYTICAL RESULTS

SESD personnel collected 24 surface soil samples and 20 subsurface soil samples (including a duplicate

sample) from various locations on the Gulf States property. One background surface soil sample (GS-01-SS)

and one background subsurface soil sample (GS-01-SB) were collected from a location northeast of the

facility. The background samples were collected from Jackson Preparatory School at 3100 Lakeland Drive.

Locations for surface and subsurface soil samples collected on-site are presented in Tables 1 and 2 of this

report, and in Reference 1, Appendix A, Figure 1.

Soil data is compared to background samples and is considered elevated if the compound is three times the

background concentration. In the case where a constituent is undetected in the background sample, any

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concentration equal to or greater than the SQL is considered to be elevated. Specific findings regarding sample results are summarized below and presented in Tables 5 through 8.

#### Surface Soil

- SVOCs were detected at elevated concentrations in eleven of the twenty-three on-site surface soil samples. Constituents detected include acenaphthylene, anthracene, benzo(a)anthracene, benzo(b)fluoranthene, benzo(g,h,i)perylene, benzo(k)fluoranthene, benzo-a-pyrene, carbazole, chrysene, dibenzo(a,h)anthracene, fluoranthene, indeno(1,2,3-cd)pyrene, phenanthrene, and pyrene. Elevated constituent concentrations ranged from 380J micrograms per kilogram (μg/kg) to 12,000J μg/kg. The majority of SVOC constituents were detected in surface soil samples GS-05-SS, GS-07-SS, GS-08-SS, GS-10-SS, GS-11-SS, GS-14-SS, GS-15-SS, GS-16-SS, GS-17-SS, GS-23-SS, and GS-24-SS which are located on the northeastern, central, and southeastern portions of the property. Areas of contamination are located where boxcars containing wood treated with creosote were stored on the railroad tracks (Refs. 1, Appendix A, Figure 1; 11).
- Endrin, endrin aldehyde, endrin ketone, and 4,4'-DDT were the only pesticides detected at elevated concentrations in the surface soil samples collected. Pesticide concentrations ranged from 6.4J μg/kg to 150 μg/kg. Pesticides were detected at elevated concentrations in eight of the twenty-three on-site surface soil samples.
- PCBs were not detected in the on-site surface soil samples.
- Methyl ethyl ketone (MEK) is the only volatile organic compounds (VOC) detected at an elevated concentration in the surface soil samples. MEK concentrations in samples GS-02-SS, GS-11-SS, GS-12-SS, GS-15-SS, GS-17-SS, GS-18-SS, GS-19-SS, GS-20-SD, GS-21-SS, and GS-23-SS ranged from 12μg/kg to 31J μg/kg. MEK is a common laboratory contaminant, and cannot be directly attributed to the processes at the facility.
- Inorganic constituents detected at elevated concentrations in on-site surface soil samples include antimony, barium, beryllium, cadmium, copper, lead, manganese, and zinc. Elevated constituent

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concentrations ranged from 0.09 milligrams per kilogram (mg/kg) to 5,500 mg/kg. Samples containing elevated levels of one or more constituent include GS-02-SS, GS-03-SS, GS-08-SS, GS-10-SS, GS-11-SS, GS-13-SS, GS-14-SS, GS-15-SS, GS-19-SS, GS-20-SS, GS-21-SS, and GS-24-SS.

#### Subsurface Soil

- Extractable SVOCs were detected at elevated concentrations in four of the twenty on-site surface soil samples. Constituents detected include acenaphthylene, anthracene, benzo(a) anthracene, benzo(b) fluoranthene, benzo(g,h,i) perylene, benzo(k) fluoranthene, benzo-a-pyrene, bis(2-ethylhexyl) phthalate, carbazole, chrysene, dibenzo(a,h) anthracene, fluoranthene, indeno(1,2,3-cd) pyrene, phenanthrene, and pyrene. Elevated constituent concentrations ranged from 420J µg/kg to 37,000 µg/kg. Elevated constituents were detected in samples GS-05-SB, GS-08-SB, GS-10-SB, GS-14-SB located along the central and southeastern portion of the property.
- The pesticides 4,4'-DDE and 4,4'-DDT were detected in subsurface soil samples GS-13-SB (6.9 µg/kg) and GS-14-SB (16J µg/kg), respectively.
- PCB-1260 was detected in subsurface soil sample GS-18-SB at 42J µg/kg.
- Acetone and MEK were the only VOCs detected at elevated concentrations ranging from 11J µg/kg to 99J µg/kg in samples GS-08-SB, GS-10-SB, GS-14-SB, and GS-15-SB. Both constituents are common laboratory contaminants that cannot be directly attributed to the processes at the facility.
- The majority of elevated inorganic constituents were detected in subsurface soil sample GS-10-SB. Barium, beryllium, cadmium, cobalt, copper, lead, magnesium, nickel, and zinc were detected at concentrations ranging from 0.87 mg/kg to 2,400J mg/kg in this sample. Elevated constituents detected in samples GS-14-SB, GS-15-SB, and GS-19-SB include lead (27 mg/kg), manganese (600 mg/kg to 1,000 mg/kg), and selenium (2.3 J mg/kg).

Several unknown and miscellaneous compounds were detected in surface and subsurface soil samples as detailed in Appendix A.

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4.2 SOURCE CONCLUSIONS

Based on the analytical results for samples collected during the PA/SI, surficial soil contamination is present

at the Gulf States property. Inorganic constituents detected at elevated concentrations include antimony,

barium, beryllium, cadmium, cobalt, copper, lead, manganese, nickel, and zinc. Organic constituents

detected at elevated levels include, but are not limited to, acenaphthylene, anthracene, benzo(a) anthracene,

benzo(b) fluoranthene, benzo(g,h,i) perylene, benzo(k) fluoranthene, benzo-a-pyrene, bis(2-ethylhexyl)

phthalate, carbazole, chrysene, dibenzo(a,h)anthracene, fluoranthene, indeno(1,2,3-cd)pyrene, phenanthrene,

and pyrene. The presence of the aforementioned polynuclear aromatic hydrocarbons (PAHs) are common

constituents in creosote wood-treating operations. Site processes involved treating wood with creosote, a

wood preservative comprised of PAHs, phenol, and cresols (Ref. 12, pp. 1, 67).

5.0 PATHWAYS

This section discusses the groundwater migration, surface water migration, soil exposure, and air migration

pathways. Additionally, this section discusses the targets associated with each pathway and draws pathway-

specific conclusions. Sampling locations and analytical results for samples collected from the specific

pathways are also discussed.

5.1 GROUNDWATER MIGRATION PATHWAY

Four groundwater samples and one background groundwater sample were collected during the PA/SI.

Groundwater sampling locations are depicted on the Sample Location Map found in Reference 1 Appendix

A and are described in Table 3. Field parameters and inorganic and organic analytical results for

groundwater samples are summarized in Tables 9, 10, and 11, respectively, located in Appendix A of this

report following Section 6.0.

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## 5.1.1 Geologic and Hydrogeologic Setting

The property is located entirely within the East Gulf Coastal Plain province of North America and, more specifically, within the Jackson Prairie Belt physiographic province (Refs. 7, p. 2; 15, p. 269; 16, p. 23). The property is located along the western border of Rankin County, northeast of Jackson, and east of the Pearl River (Ref. 3). The topography of Rankin County ranges from gently rolling to steep with elevations ranging from 612 feet above mean sea level (amsl) to 220 feet amsl, the property is located at approximately 265 feet amsl (Refs. 3; 7, p. 3). The soil that directly underlies the property is comprised of the Cascilla-Arkabutla soil group which is classified as nearly level, well-drained and somewhat poorly drained, silty soil. These soils typically occur along the flood plains of the Pearl River and its tributaries. The nearly linear flood plain surface within the vicinity of the property is irregularly broken by old river runs, natural levees, sloughs, chutes, and scarps. The slope of the soils ranges from 0 to 2 percent and the average thickness of the soil is approximately 40 feet (Refs. 7, p. 8; 17, p. 32). The property is underlain in descending stratigraphic order by the Cockfield Formation, the Cook Mountain Formation, the Kosciusko Formation (also known as the Sparta Sand), the Zilpha and Winona Formations, the Tallahatta formation, and the Wilcox Group (Ref. 18).

The Claiborne Group consists of the Cockfield Formation, the Cook Mountain Formation, the Kosciusko Formation, the Zilpha and Winona Formations, and the Tallahatta formation (Ref. 18). The Cockfield Formation consists of irregularly bedded laminated lignitic clay, sand, and lignite that is slightly glauconitic (Ref. 18). Based on drillers' logs, the top of the Cockfield is located at approximately 40 feet below ground surface (bgs) and is approximately 133 feet thick in the vicinity of the property (Ref. 19). The Cook Mountain Formation underlies the Cockfield and consists of marl, limestone, glauconitic sand, and chocolate colored clay (Ref. 18). The Sparta Sand (Kosciusko Formation) consists of an irregularly bedded sand with clay and some quartzite. The Sparta Sand is approximately 300 feet thick in the area, but can reach a thickness in excess of 800 feet in southwestern Hinds County (Ref. 16, p. 15). The Zilpha and Winona Formations underlie the Sparta Sand and consist of a chocolate colored clay that contains glauconitic sand and a highly glauconitic more or less clayey sand, respectively (Ref. 18). The Zilpha Formation ranges in thickness from 200 feet on the Jackson Dome to 420 feet in the southwestern portion of Rankin County. The Winona Formation ranges from 10 to 15 feet thick over the Jackson Dome to up to 65 feet thick in other

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portions of the county (Ref. 16, p. 190). The Tallahatta Formation underlies the Winona Formation and consists of predominately glauconitic claystone and clay with lenses of sand and some sandstone (Ref. 18). The Wilcox Group underlies the Claiborne Group and consists of irregularly bedded fine to coarse sand, more or less lignitic clay and lignite (Ref. 18). The Wilcox Group varies in thickness from 1,100 feet to 1,300 feet over the Jackson Dome, and attains a thickness of 2,830 feet in southern Rankin County (Ref. 16, p. 188).

Three aquifers are available for the development of moderate to large groundwater supplies in the vicinity of the property. The aquifers in descending stratigraphic order, are the Cockfield Formation, the Sparta Sand, and the Wilcox Group. All of these aquifers are part of the Eocene aquifer system in Mississippi and extend to the west, southwest, and south, and contain freshwater in approximately 50 percent of the State. All of the aquifers are regional in extent and all except the Cockfield and the lower Wilcox merge northward into a single aquifer south of Memphis, Tennessee (Ref. 15, p. 274). The formations dip southwest at approximately 15 to 25 feet per mile toward the Mississippi Embayment and the Mississippi River, and the groundwater flow generally follows this regional trend (Ref. 17, p. 4). Within the geologic column, the water-bearing sand beds are interbedded with shale of both marine and continental origin, fossiliferous limestone, and calcareous sandstone. Strata that were deposited by marine origin generally consist of clay, and they form aquicludes (or confining layers) between the water-bearing sands. The aquicludes are widespread and more uniform in thickness than the aquifers (Ref. 17, p. 4).

The Cockfield Formation is the source of more than half of the municipal water supply in the area, mainly because it is the most shallow aquifer in the area (Ref. 17, pp. 1, 32). The top of the formation is located beneath the alluvial soil deposits at approximately 40 feet bgs in the vicinity of the property, ranges from 80 to 140 feet thick, and is unconfined (Refs. 17, p. 32; 19). Municipal water supplies for several small towns are obtained from the Cockfield, some wells yielding as much as 500 gallons per minute (gpm) (Ref. 17, p. 1). The hydraulic conductivity of the Cockfield is approximately  $1 \times 10^{-2}$  centimeters per second (cm/s) (Ref. 20, p. 29). The Cook Mountain Formation underlies the Cockfield aquifer and consists of marl, limestone, glauconitic sand, and chocolate colored clay (Ref. 18). The Cook Mountain Formation was deposited in a marine environment, exhibits a high clay content, and serves as a confining layer between the overlying

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Cockfield Formation and the underlying Sparta Sand (Ref. 17, p. 4).

The Sparta Sand underlies the Cook Mountain Formation in the vicinity of the property and is approximately

300 feet thick in the area, but can reach a thickness in excess of 800 feet in southwestern Hinds County. The

Sparta Sand is the most intensively developed aquifer in the vicinity of the property. The Sparta exhibits a

hydraulic conductivity of approximately 1x10<sup>-2</sup> cm/s (Refs. 17, pp. 15, 32; 20, p. 29).

The Zilpha and Winona Formations underlie the Sparta Sand and consist of a chocolate colored clay that

contains glauconitic sand and a highly glauconitic more or less clayey sand, respectively (Ref. 18). Due to

the high clay content and marine origin of these formations, they serve as the lower confining layer for the

Sparta Sand and the upper confining layer for the Wilcox Group in the vicinity of the property. These

formations underlie the area at depths that range from 650 feet in northeastern Madison County to 2,600 feet

in southwestern Hinds County. The depth of the formations throughout the remainder of the study area

ranges from 420 to 570 feet (Ref. 17, p. 15).

The Wilcox Group contains a large reserve of soft water that has been tapped by only a few small-supply

wells. The water is more highly mineralized and is warmer than that in the more shallow aquifers. The water

in the aquifer is of good quality in Madison and northern Rankin Counties, but the quality deteriorates down

the dip in Hinds County (Ref. 17, p. 1). In counties to the northeast, the results of several pumping tests

indicate that the sands in the Wilcox aquifer are probably as permeable as the Sparta Sand (Ref. 17, p. 15).

5.1.2 Groundwater Sampling Locations and Analytical Results

SESD personnel collected four groundwater samples from temporary monitoring wells installed on the Gulf

States property. The background groundwater sample was collected from the Jackson Preparatory School

at 3100 Lakeland Drive located northeast of the facility. Locations for the groundwater samples are detailed

in Table 3 and depicted in Reference 1, Appendix A, Figure 1. Within this report, Appendix A, Table 9 lists

the field parameters for the groundwater samples. Sample results are summarized in Tables 10 and 11

located in Appendix A of this report.

Inorganic constituents detected at elevated concentrations when compared to background concentrations

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include aluminum, barium, cobalt, magnesium, manganese, vanadium, and zinc. Concentrations of

inorganic constituents ranged from 1.9 micrograms per liter (µg/L) to 8,200 µg/L. Barium, cobalt,

magnesium, manganese, and zinc were also detected at elevated concentrations in on-site soil samples

and can therefore be attributed to the site.

• MEK is the only organic compound detected at an elevated concentration in GS-04-GW at 7.2 J μg/L.

• No SVOCs, pesticides, or PCBs were detected in any of the groundwater samples:

5.1.3 Groundwater Targets

During the June 1993 Phase I investigation at the property, substances detected in groundwater samples

collected from a temporary monitoring well included 2,4-dimethylphenol, naphthalene, dibenzofuran,

fluorene, phenanthrene, fluoranthene, and pyrene at concentrations ranging from 79 micrograms per liter

 $(\mu g/L)$  to 1,279  $\mu g/L$  (Ref. 4, p. 26).

The majority of local residents obtain drinking water from the City of Flowood Water Department (CFWD).

The CFWD obtains its drinking water from six wells screened in the Cockfield Formation and the Sparta

Sand aquifers, which have depths ranging from 562 feet to 1,185 feet. Water from the wells is blended prior

to distribution within the system. CFWD serves approximately 5,400 connections (Ref. 21). According to

the 2000 Census, there are 2.62 people per household in Rankin County (Ref. 22). Based on the number of

service connections and the number of people per household, CFWD serves approximately 14,148 people

(2.62 people per household x 5,400 connections with an average of 2,358 people served per well) (Refs. 21;

22). The CFWD well distribution is as follows: 0 to 0.25 mile: 0 wells; 0.25 to 0.50 mile: 0 wells; 0.50 to

1 mile: 1 well; 1 to 2 miles: 0 wells; 2 to 3 miles: 1 well; 3 to 4 miles: 1 well. Therefore, the CFWD

population distribution is as follows: 0 to 0.25 mile, 0 persons; 0.25 to 0.50 mile, 0 persons; 0.50 to 1 mile,

2,358 persons; 1 to 2 miles, 0 persons; 2 to 3 miles, 2,358 persons; 3 to 4 miles, 2,358 persons (Refs. 3; 21;

22). No private wells were identified in the surrounding neighborhoods during the June 2002 site

reconnaissance (Ref. 5, p. 11).

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5.1.4 Groundwater Conclusions

The only site-attributable constituents detected in the groundwater samples collected from the surficial

aquifer during the PA/SI were barium, cobalt, manganese, and zinc. The majority of people in the area

receive their drinking water from the City of Flowood Water Department. CFWD's wells are screened in

the Cockfield Formation and the Sparta aquifers below a confining layer. Due to the small number of people

served water by the CFWD and the fact that the municipal wells are screened at depths greater than 550 feet,

the groundwater pathway does not appear to be a viable migration route.

5.2 SURFACE WATER MIGRATION PATHWAY

Six sediment samples were collected during the PA/SI to document the migration of on-site contaminants

into the marsh located west of the Gulf States property, and the Pearl River (Ref. 3).

5.2.1 Hydrologic Setting

Currently, a levee separates the Gulf States property from the marsh area located west of the property (Refs.

3; 5, pp. 7, 8). The construction date of the levee is unknown, however, the levee was in existence in 1955

(Ref. 4, p. vii). The surface water migration pathway prior to the construction of the levee, appears to have

flowed west toward the marsh and the tributaries which flow into the Pearl River (Ref. 3). There is a

drainage ditch on the western portion of the property which flows north to south. The drainage ditch is

located to the east of the levee (Ref. 5, pp. 7, 8). Currently, it appears that the on-site drainage flows into

this drainage ditch. Prior to the construction of the levee, it appears that creosote was discharged to the

marsh areas west of the property, and these areas have been defined as creosote slough by the U.S.

Geological Survey (Refs. 3; 5, p. 4). During the operation of the facility, the Pearl River was re-channeled

to flow further to the west of the facility and the crossote slough areas (Refs. 3; 11). The surface water

pathway includes the marshland and tributaries which flow approximately 0.8 mile to the Pearl River. The

15-mile downstream surface water pathway is completed in the Pearl River which flows at an average rate

of 759 cubic feet per second (Refs. 3;23). The property itself is within a flood area that is undefined, perhaps

due to the construction of the levee. However, the marsh area adjacent to the facility is within the 100-year

flood plain (Ref. 24). A City of Jackson surface water intake is located approximately 0.75-mile northwest

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of the Gulf States property on the Pearl River (Ref. 5, p. 4).

5.2.2 Sediment Sampling Locations and Analytical Results

SESD personnel collected five sediment samples from the marsh adjacent to the Gulf States property and the

Pearl River. Two background sediment samples (one from the Pearl River and one from a marsh) and one

control sediment sample were collected for comparison to the downstream samples. The background and

control sediment sample locations are detailed in Table 4 and depicted in Reference 1, Appendix A, Figure

1. Specific findings regarding sample results are summarized below and presented in Tables 12 through 13.

• Several SVOCs were detected at elevated concentrations in the sediment samples. Such constituents

include acenaphthene, acenaphthylene, anthracene, benzo(a) anthracene, benzo(b) fluoranthene,

benzo(k) fluoranthene, benzo-a-pyrene, chrysene, dibenzo(a,h) anthracene, fluoranthene, fluorene, 2-

methylnaphthalene, naphthalene, phenanthrene, and pyrene. Constituent concentrations ranged from

2 μg/kg to 560J μg/kg. Elevated constituents were detected in all of the sediment samples collected

from the marsh adjacent to the Gulf States property.

VOCs including acetone, carbon disulfide, and MEK were detected at elevated concentrations in

sediment samples GS-03-SD and GS-04-SD.

• The pesticides 4,4'-DDE (7.9 μg/kg) and beta-BHC (12 μg/kg) were detected in sediment sample GS-03-

SD and GS-04-SD, respectively.

Inorganic constituents detected at elevated concentrations in sediment samples are barium, beryllium,

lead, and zinc. All four constituents were detected in sample GS-03-SD at concentrations ranging from

1.3 mg/kg to 300 mg/kg. An elevated level of beryllium was also detected in sample GS-04-SD (1.2

mg/kg).

Several unknown and miscellaneous compounds were detected in the sediment samples as detailed in

Appendix A of this report.

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5.2.3 Surface Water Targets

Federally endangered or threatened species within the Pearl River are the Ringed map turtle (Graptemys

oculifera) and the Gopher tortoise (Gopherus polyphemus) (Refs. 25, p. 2; 26, pp. 20, 21). Although their

exact location is unknown, other endangered/threatened species potentially within the area include the Bayou

darter (Etheostoma rubrum); Pallid sturgeon (Scaphirhynchus albus); Southern combshell (Epioblasma

penita) (Ref. 26, pp. 23, 30, 31). Approximately 2.5 miles of wetland frontage is present along the Pearl

River (Ref. 3). Recreational fishing occurs in both the marsh and Pearl River. Commercial fishing also

occurs along the Pearl River (Ref. 5, pp. 3, 4). There is a fish tissue advisory in effect for the Pearl River

from Highway 25 near Carthage, downstream to the Leake County Water Park. The advisory recommends

limited consumption of largemouth bass and large catfish due to mercury contamination (Ref. 27).

**5.2.4 Surface Water Conclusions** 

Sediment samples collected from the marsh located west of the Gulf States property revealed elevated levels

of site-attributable contaminants. The constituents include acenaphthene, acenaphthylene, anthracene,

benzo(a) anthracene, benzo(b) fluoranthene, benzo(k) fluoranthene, benzo-a-pyrene, chrysene, dibenzo(a,h)

anthracene, fluoranthene, fluorene, 2-methylnaphthalene, naphthalene, phenanthrene, and pyrene at

concentrations ranging from 2 µg/kg to 560J µg/kg. These constituents are common constituents in the

composition of creosote and are commonly found at wood treating facilities. Such constituents pose a threat

to the marsh and the Pearl River due to the recreational fishing that occurs in these water bodies. Due to the

detection of the aforementioned constituents in the nearby surface water bodies, the surface water pathway

is a viable migration route for contaminants.

5.3 SOIL EXPOSURE PATHWAY AND AIR MIGRATION PATHWAY

Twenty-three surface soil and 20 subsurface soil samples were collected during the PA/SI at the Gulf States

property. Surface and subsurface soil sample results were discussed in Section 4.0.

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5.3.1 Physical Conditions

The 141-acre Gulf States property is zoned for heavy industrial use (I-2) (Refs. 5, p. 2; Ref. 13, p. 1). The

property currently consists of a horse farm and a company that sells reinforcing steel and wire mesh (Ref.

5, pp. i, 1). The facility is bound by railroad tracks to the north and east, an adjacent business to the south,

and marshland/tributary to the Pearl River to the west (Ref. 3; 11). The facility is bound by natural barriers

and portions of the facility property are fenced. The fenced portions of the property include the area

surrounding the ConSteel building, and the horse grazing area (Ref. 5, p. 8). Access to the entire property

is unrestricted.

During the 1993 Phase I conducted at the property, seven soil borings were constructed ranging in depth from

0 to 8 feet bgs. Of the seven borings, five samples were collected and analyzed for PAHs. Constituents

detected in the soil samples included napthalene, 2-methylnapthalene, dibenzofuran, fluorene, phenanthrene,

and fluoranthene. Organic concentrations ranged from below the detection limit to 604 (mg/kg) (Ref. 4).

**5.3.2** Sampling Locations and Analytical Results

Surface and subsurface soil sampling locations and analytical results are discussed under Source Sampling

in Section 4.1. No air samples were collected at the facility.

5.3.3 Soil and Air Targets

The facility currently houses a horse farm and the ConSteel Company (Ref. 5, p. 1). The nearest residences

are located east of the property across Highway 15 (Ref. 3; 4). There are no schools, daycare facilities, or

residences within 200 feet of the property (Refs. 3; 5, p. 11). However, the two businesses on the property

are operational and the potentially exposed population includes the workers at these two facilities (Ref. 5,

p.1).

According to a LandView<sup>®</sup> Census Data report and a house count, approximately 805 persons reside within

4 radial miles of the Gulf States property. The estimated population distribution within 4 radial miles of the

Gulf States property is as follows: 0 to 0.25 mile, 118 persons; 0.25 to 0.50 mile, 45 persons; 0.50 to 1 mile,

152 persons; 1 to 2 miles, 87 persons; 2 to 3 miles, 44 persons; and 3 to 4 miles, 359 persons (Ref. 28; 29).

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Several federal threatened and endangered species are suspected within the vicinity of the Gulf States

property. Such species may include the Indiana bat (Myotis sodalis). However, their exact locations have

not been identified (Ref. 26, p. 2). Approximately 410 acres of wetlands are located within 4 miles of the

Gulf States property. The wetlands distribution is as follows: 0 to 0.25 mile, 10 acres; 0.25 to 0.50 mile, 15

acres; 0.50 to 1 mile, 185 acres; 1 to 2 miles, 190 acres; 2 to 3 miles, 10 acres; 3 to 4 miles, 0 acres (Ref. 3).

5.3.4 Soil and Air Conclusions

Elevated constituents were detected in the surface soil samples collected from the property; however, the

property is not residential, and has been zoned heavy industrial. Based on the nature of the constituents

detected at the property, the potential exposure to employees and local residences via soil or air is considered

minimal.

6.0 SUMMARY AND CONCLUSIONS

The former 141-acre Gulf States Creosoting Company is located at 1625 Flowood Drive (Mississippi Hwy

468), Flowood, Rankin County, Mississippi. The facility is currently operating as a horse farm and ConSteel.

Both current operations appear to be located on portions of the property where creosoting operations were

conducted. Currently, the southernmost building on the property is owned by ConSteel. The northernmost

structures (i.e. metal barn and shed) are used to house tractors and helicopters for a local news station, and

are part of the horse farm. The horse farm has been operating on the property for approximately 7 years.

Gulf States owned the property as early as 1929 and operated as a wood treating facility until the mid 1950s.

In July 1958, American Creosoting Corporation obtained portions of the property and operated there until

the late 1950s. In June 1959, W.G. Avery Body Company purchased portions of the property and operated

a body shop on the site. There are several tracts of land on the Gulf States property and some are currently

occupied. The horse farm tract is owned and operated by Mr. Jim Webb; Webb owns the portion of the

property just north of the main entrance to the property. ConSteel owns a portion of the property due south

of the main entrance. ConSteel appears to be operating on the portion of the property that it owns. The

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onsite levee is owned by the Levee Board - which is a joint venture between the state and Rankin County.

Mr. John McGowan owns the marsh area due west of the levee. McGowan has plans of developing the

marsh areas as residential subdivisions.

During Gulf States operations, railroad cross ties were treated at the facility and transported on and off-site

by means of railroad box cars. Creosote was applied to the wood by commercial pressure treatment or by

home and farm dipping. Creosote is a wood preservative used to treat railroad ties, telephone poles, marine

pilings, and fence posts.

Based on the analytical results for the samples collected during the PA/SI, soil contamination is present at

the Gulf States property. Inorganic constituents detected at elevated concentrations include antimony,

barium, beryllium, cadmium, cobalt, copper, lead, manganese, nickel, and zinc. Organic constituents

detected at elevated levels include, but are not limited to, acenaphthylene, anthracene, benzo(a)anthracene,

benzo(b)fluoranthene, benzo(g,h,i)perylene, benzo(k)fluoranthene, benzo-a-pyrene, bis(2-

ethylhexyl)phthalate, carbazole, chrysene, dibenzo(a,h)anthracene, fluoranthene, indeno(1,2,3-cd) pyrene,

phenanthrene, and pyrene. The presence of the aforementioned PAHs are attributable to the site operations.

The only site-attributable constituents detected in the surficial aquifer were barium, cobalt, manganese, and

zinc. The majority of people in the area receive their drinking water from the City of Flowood Water

Department. CFWD's wells are screened in the Cockfield Formation and the Sparta aquifers beneath a

substantial a confining layer. Due to the small number of people served water by the CFWD and the fact

that the municipal wells are screened at depths greater than 550 feet, groundwater does not appear to be a

viable migration route.

Sediment samples collected from the marsh located west of the Gulf States property exhibited elevated levels

of constituents detected in on-site soils and are common constituents used in the wood preserving industry.

Such constituents include acenaphthene, acenaphthylene, anthracene, benzo(a) anthracene, benzo(b)

fluoranthene, benzo(k) fluoranthene, benzo-a-pyrene, chrysene, dibenzo(a,h) anthracene, fluoranthene,

fluorene, 2-methylnaphthalene, naphthalene, phenanthrene, and pyrene at concentrations ranging from 2

μg/kg to 560J μg/kg. Such constituents pose a threat to both the marsh and the Pearl River because of the

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recreational fishing that occurs on these water bodies. Due to the detection of the aforementioned constituents, the surface water pathway is a viable migration route for contamination.

Surface soil contamination is present at the property. However, because the property is primarily industrial and not residential, and the non-volatile nature of the contamination, the exposure to employees and local residences by means of soil or air is considered to be minimal.

Based on the analytical results for the samples collected during the PA/SI, <u>further CERCLA action is</u> recommended for the Gulf States property.

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## APPENDIX A

## ANALYTICAL DATA TABLES

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## APPENDIX A

## ANALYTICAL DATA TABLES

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## TABLE 5 SUMMARY OF INORGANIC ANALYTICAL RESULTS SURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

	ĺ	⊮Background∄	345, 25-15	57.58 (ST			Transfer		On-site Sar	nples	A39100年5月27	are the second	ecumum, dat. 3	THE PLANT OF THE
COMPOUND	UNITS	GS01SS	GS02SS	GS03SS	GS04SS	GS05SS	GS06SS		GS08SS	GS09SS	GS10SS	GS11SS	GS12SS	GS13SS
METALS *	12	海绵红 编辑统计		种类类的	ry baar	医松叶与中的	和60年年12日	的的技术的经	No. of the		ere de la company		1774	2007年7月44年74
ALUMINUM	MG/KG	4700 J	7800 J	8900 J	4300 J	3100 J	6700 J	2800 J	1800 J	2500	4700 J	4400 J	5500 J	4400 J
ANTIMONY	MG/KG	0.56 UJ	0.64 R			•				1.2 R	1 R			0.6 R
ARSENIC	MG/KG	6.3	3.3 J	11	2	13	2	3.6	3.4	0.95 R	6.8	3.7	4.4	6.3
BARIUM	MG/KG	72	120	67 _	60	37	85	34	21	25	120	54	83	58
BERYLLIUM	MG/KG	0.33	0.94	0.94	0.37	0.3	0.28	0.2	0.19	0.15	0.46	0.44	0.46	0.54
CADMIUM	MG/KG	0.05 U		-		-	<b></b> -		<b>00</b>		0.34	0.19		0.16 R
CALCIUM	MG/KG	1100	690	790	690	500	1200	810	800	340	2600	660	760	1200
CHROMIUM	MG/KG	8.4	6.1	20	6	7.5	7.7	4.5	4	4.1	23	12	8.9	7.8
COBALT	MG/KG	5.8	17	8.1	1.7	6.6	1.4	2	1.5	0.95	4.8	5.4	7.3	4.2
COPPER	MG/KG	6	2 J		<b></b> .			-			3962-41975			7.2 J
IRON	MG/KG	10000	8400	27000 J	4500 J	15000 J	5600 J	5400 J	4700 J	2800	14000	11000	11000 J	12000
LEAD	MG/KG	11 J	17	120	18	19	17	11	12	8.8	少划 #80 ##	290	25	<b>1</b> 2.2.5 <b>33</b> .
MAGNESIUM	MG/KG	520 J	470 J	680	280	180	550	250	180	160 J	780 J	260 J	380	280 J
MANGANESE	MG/KG	630	2800	550	140	610	170	180	230	28	500	880	990	940
POTASSIUM	MG/KG	380 J	240 J	350	200	120	430	170	150	120 J	560 J	330 J	280	390 J
SELENIUM	MG/KG	1.3	0.93 R	1.9 J	'				••		0.98 R	0.77 R		1.2
SILVER	MG/KG	0.39 R	0.71 R	0.63	0.26 R	0.43	0.29		<i></i>	0.28 R	0.49	0.38	0.43	0.33 R
SODUIM	MG/KG	270	300_	82 J	120 J	58 J	110	63		300	320	200	91	230
VANADIUM	MG/KG	17	15	46	11	18	. 13	8.1	7.1	6.6	18	16	22	17
ZINC	MG/KG	29	47	34	27	23	35	24	26	19	160	840	28	72

### Notes:

GS - Gulf States Creosoting

SS - surface soil sample

MG/KG - milligrams per kilogram .

U - value is below the reporting limit

J - estimated value

R - data is unusable

shading - elevated concentrations of constituents

## TABLE 5 (CONTINUED) SUMMARY OF INORGANIC ANALYTICAL RESULTS SURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

		Background	6 (2 a) ( ( a) ( )	lowers.	STAINING.	SERVICE AND	grown	્રી ં્રા	n-site Sam	oles 🔆 🔆		in, C. S. Marce S	Y Julius In 1911.	
COMPOUND	UNITS	GS01SS	GS14SS	GS15SS	GS16SS	GS17SS	GS18SS	GS19SS	GS20SS	GS20SD*	GS21SS	GS22SS	GS23SS	GS24SS
METALS	Silver - S.		T. 120	學生學的 斯瑟			20 C * 3 m S				《北美公徽》(	Huding Lea	<b>聚</b> 的原料。这	Contain Party
ALUMINUM	MG/KG	4700 J	4800 J	8000 J	7700 J	9000 J	4900 J	8500 J	6700 J	6100 J	6800 J	4200 J	7500 J	4900 J
ANTIMONY	MG/KG	0.56 UJ	0.69 J			••					7-			
ARSENIC	MG/KG	6.3	9	6.9	5.5 J	5.5	5.4	6.5	5.8	4.5	2.7	4	5	6.5
BARIUM	MG/KG	72	65	72	50	120	53	89	410	130	88	42	83	120
BERYLLIUM	MG/KG	0.33	0.44	0.71	0.43	0.94	0.35	0.73	1.2	0.76	0.81	0.43	0.5	0.61
CADMIUM	MG/KG	0.05 U	0.11 R	<i>≫</i> 0.1		0.07 R		-	0.17				0.1 R	0.16
CALCIUM	MG/KG	1100	1800	2000	1200	2300	890	2000	1700	1400	190	280	960	1600
CHROMIUM	MG/KG	8.4	8.6	11	10	8	. 9.9	8.3	8.2	7.2	5.6	9.4	12	12
COBALT	MG/KG	5.8	7.3	11	4	12	7.6	13	17	6.8	14	5.5	6.8	6.8
COPPER	MG/KG	6	·					-					12 J	
RON	MG/KG	10000	11000	12000	12000	12000	10000	12000	12000	8800	7300	9700 J	13000	17000 J
LEAD	MG/KG	11 J	24	26	15 J	20 J	15 J	19 J	24 J	22 J	13 J	9.8	24	. 28
MAGNESIUM	MG/KG	520 J	340 J	430 J	510 J	490 J	340 J	530 J	380 J	410 J	270 J	220	630 J	450
MANGANESE	MG/KG	630	870	1800	240	2500	590	1900	5500	1800	2200	470	920	890
POTASSIUM _	MG/KG	380 J	200 J	580_J	490 J	620 J	180 J	550 J	240 J	250 J	250 J	170	480 J	370
SELENIUM	MG/KG	1.3	0.63 R	1.1 R	1.3	1.5 J	1.1	1.2	1.3 J	1.1 R	0.8		1.1 R	
SILVER	MG/KG	0.39 R	0.37	0.5	0.39 R	0.51 R	0.29	0.53 R	0.9 R	0.42	0.52	0.2	0.46 R	0.47
SODUIM	MG/KG	270	300	310	270	280	310	290	290	290	300		340	86 J
VANADIUM	MG/KG	<u>1</u> 7	21	23	22	23	20	-22	22	17	13	21	23	. 26
ZINC	MG/KG	29	44	48	33	65	34	52 .	. 88	63	24	19	49	<b>89</b>

### Notes:

GS - Gulf States Creosoting

SS - surface soil sample

MG/KG - milligrams per kilogram

U - value is below the reporting limit

J - estimated value

R - data is unusable

shading - elevated concentrations of constituents

- Sample GS20SD is a duplicate of sample GS20SS.

## TABLE 6 SUMMARY OF ORGANIC ANALYTICAL RESULTS SURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

•	1									•		•		
		Background	Company of	translation of	Kardakimi beli	البهر مها معصلي	Lathery Lynn		::::On-site 8ei	nples (2005)		医基金性线 医线点	A STATE OF THE	તો, ૧૯૧૬ ભોજરીફેફ ફ
		ì									T		ĺ	
COMPOUND	UNITS	G501SS	GS02SS	GS03S8	G\$04\$\$	GS0588	QS0688	GS07SS	GS08SS	GS09SS	GS10SS	GS11SS	GS12SS	GS13SS
VOLATRES	and the many the Court of the C	a file stage of the original		340.007, provi	William Thomps		POW WELL	50 144 - 1136	- 25 Table 17 (9)	CHESTOLIC CONTRACT	17.27× 28601 18.28	kapangalah tilabak i	2. 400 Oct 4000 A	ACCULEMENT
ACETONE	UG/KG	100 J	180 J	25 J	28 J	91 J	82 J	87 J	74 J		84 J	220 J	T	200 J
METHYL ACETATE	UG/KG	11 U			-		<del></del>	-	-	-	T	_	<u> </u>	
METHYL ETHYL KETONE	Jug/kg	11 U	a 21.640053				L=	L =	I	L	I	55-55821 J.S.	2-45-414 J	15 J
MISCELLANEOUS VOLATILE ORGANICS	Const. St. Company and the second second	a rango ( ) ki pigamenej	46, 54, 5 . 35 . 38 . 38 ·	on Christian 18	addition of the San	ar a landina di inagi	Primari U.A.	Carles and P	wind in the No. 1 of 4.632	. و يا خون پرې د مړۍ د ويوان د	and the state of the state of	Car. Printing of the	1 4 M 1- 132 Merel	Ger Joseph Grand Story
UNKNOWN COMPOUNDS/NO.	IUG/KG	19 J/2	51 J/2	6 J/1			20 J/2			1	29 J/3		30 J/3	13 J/1
ACETALDEHYDE	UG/KG								12.23	<del> </del>	+	<del>                                     </del>	<del>                                     </del>	16 NJ
BUTANAL	UG/KG						1				<del></del>		5 NJ	
BUTANAL 3-METHYL-	UG/KG	f									<del> </del>			
HEXANAL	UG/KG	i			<del></del>		64 NJ				45 NJ	37 J	18 NJ	
PENTANAL	UG/KG							<del></del>	1		<b> </b>			
PROPANAL, 2-METHYL-	UG/KG								<u> </u>		Ì			
											_			
EXTRACTABLES AND														
(3-AND/OR 4-)METHYLPHENOL	UG/KG	380 U		-	<u> </u>				<u> </u>		<del> </del>	<u> </u>		
ACENAPHTHENE	UG/KG	380 U	-		-		-	170 J			97 J	<u></u>		
ACENAPHTHYLENE	UG/KG	380 U	ا ا	-		120 J	-	320 J	480 J		1.00 ×c	330 J		1000 J
ANTHRACENE	UG/KG	380 U		-	<u> </u>	470 X		50 970 m			-8:34700 ast	930 a 174		1600
BENZALDEHYDE	UG/KG	380 U								<del></del>			<u> </u>	
BENZO(A)ANTHRACENE	UG/KG	380 U			L <del>-</del>	50.770 S (A)	÷	2500	1500		::55%.2800.ide:	565 9 <b>1000</b> 307 A		4300
BENZO(B)FLUORANTHENE	UG/KG	. 380 U		-		1800 🕽 🚐	<u> </u>	4600 Jo			5200 234	2500		7300 J
BENZO(GHI)PERYLENE	UG/KG	380 U	_=			≆380 J∴	_=	«690∴	* 5800 : /#	:		3 4 5 480 to		2000
BENZO(K)FLUORANTHENE	UG/KG	380 U		-		2100 J		-5200 J.√			2300			5400 J
BENZO-A-PYRENE	UG/KG	380 U		-		G <b>610</b> ∜%≪		2100 Tes	3700∠√€⊴			800	-	3100
1,1-BIPHENYL	UG/KG		-	-				45 J						
CARBAZOLE	UG/KG	380 U		-		200 J		510 10			21. 2 480 ( **	210 J		540 J
CHRYSENE	UG/KG UG/KG	380 U	-	-		1500 Acres		. 4900 CA		<del>` -</del>		ASS 15001 10 D		6200
DIBENZO(A,H)ANTHRACENE DIBENZOFURAN	UGAKG	380 U			<del>-</del>	180 3		240 J	#1700 P. P. C.		4.5° 39 <b>750</b> (\$7	310 J	<u></u>	1100
FLUORANTHENE	UGAKG	380 U	<del>  </del>		<del></del>	2800		18200 4 (c)	≥2600 // 3.55//		120 J	140 J		11000
FLUORENE	UG/KG	380 U	<del></del>	<del></del>		2001 5:4-17	-	180 J	# 2000 / A 52/1		4700	* 47 1900 A. C.		
INDENO(1,2,3-CD)PYRENE	UG/KG	380 U				+ <b>610</b> 72561		÷1100.5	7/ <b>5000</b> (1/6/44)					
2-METHYLNAPHTHALENE	UG/KG	380 U				7010-098		120 J			59 \$ 2300 V	Kirne j 940 kir. (19		3100
NAPHTHALENE	UG/KG	380 U			<del></del> -			87 J		-	69 J	70 J		
PHENANTHRENE	UG/KG	380 U			<del>-</del>	320 J	-		563 <b>010.</b> 30.333		110 J	473014 <b>840</b> 33703		1600
PENTACHLOROPHENOL	UG/KG	970 LU			<del></del> -	320 J		# 42LU -242	76 GIO 33 223			640 J		
PYRENE	UG/KG	380 U	-	- 1	<del></del> _	1500 (2.4)	=	> 8200 P.T.			690 J	640 J	<del>- :-</del>	6800
THENE	Journa	360 0			<del></del>	v (000 (23×4)		P OE OO STORY	#2000 (ASH Q)		in Literature	(490,01700;93%)		8800
MISCELLANEOUS EXTRACTABLES	WW Westernamer Co	our Ting displaying	orthographic	ed all a	Mass: last) (June	888 X 60 5 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	GARREST W. SAME	N. 411 (42)	<b>2004年,1907年</b> 。	configuration be	en het Galderide	AND PARTY OF THE P	(PO) 0.28年中17月	AND WAR SHOW
UNKNOWN COMPOUNDS/NO .	UG/KG	1400 J/12	2200 J/5	91 J/1	2000 J/7	2700 J/11	2800 J/14	5900 J/8	23000 J/17	1900 J/10		3600 J/12	410 J/3	19000 J/12
STIGMAST-4-EN-3-ONE	UG/KG		190 NJ		350 J		460 NJ							
UNKOWN ALCOHOL	UG/KG				92 J						1			
UNKNOWN KETONES/NO.	UG/KG				250 · J/2									
METHYLANTHRACENE	UG/KG					190 J					<u> </u>			
9,10-ANTHRACENEDIONE	UG/KG					260 NJ		890 NJ			660 NJ	260 NJ		890 NJ
PHENANTHRENE, 1-METHYL	UG/KG													1100 NJ
PHENANTHRENE, 2-METHYL	UG/KG													
PHENANTHRENE, 2,3-DIMETHYL	UG/KG		1			160 NJ		870 NJ				-		1300 NJ
PHENANTHRENE, 4,5-DIMETHYL	UG/KG							830 NJ	1					
ANTHRACENE, 2-METHYL	UG/KG							1000 J						
BENZOFLUORENE	UG/KG					160 J			. 920 J					
I 1 H-BENZOIAIFLUORENE	UG/KG					470 NJ		740 NJ			2200 NJ			
11 H-BENZO(B)FLUORENE	UG/KG													1100 NJ

## TABLE 6 (CONTINUED) SUMMARY OF ORGANIC ANALYTICAL RESULTS SURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

		Beckground	<b>本</b> 等。可以2.45	Mark States	ing true professions	2 m. 1987 to 20 pict	Maria (Albania)	A 147 ( 15	On alte Ser	mples were seen	er Kildir inderktore i	ري. دون د دون ايا د دون وي	PARTITE SEE	The contraction of
COMPOUND	UNITS	G\$0155	GS025S	GS03S6		GS058S				GS09SS	G\$1088	GS11SS	GS12SS	GS13S8
2,2'-BINAPHTHALENE	UG/KG							· ·			560 N	†···	<del> </del>	560 NJ
1,2:3,4-DIBENZPYRENE	UG/KG				1	<del>                                     </del>		i	<del>   </del>	1	7——	<del> </del>		
1,2:4,5-DIBENZPYRENE	UG/KG					i		<del> </del>	880 NJ	1				
1-DOCOSENE	UG/KG	<del> </del>				·		<del> </del>				1		<del>                                     </del>
1-EICOSANOL	· UG/KG	1	1											
PENTADECANOIC ACID	UG/KG				· · · · ·					97 NJ	1			1
PERYLENE	UG/KG								3900 NJ		1000 NJ			
PYRENE, 1-METHYL	UG/KG	·							880 NJ		710 NJ	390 NJ		1400 NJ
PYRENE, 2-METHYL	UG/KG					250 NJ		1200 NJ	1000 NJ		1700 NJ			660 NJ
PYRENE, 1,3-DIMETHYL	UG/KG					280 NJ		1992 119						
PYRENE, 4-METHYL	UG/KG		· —			200 110		740 NJ		<del></del>	<del> </del>			
BENZIAIANTHRACENE, 7-METHYL-	UG/KG							7-10 110		<del></del>	+			t —
BENZ(A)ANTHRACENE-7, 12-DIONE	UG/KG							540 NJ			+	330 NJ		
BENZOBINAPHTHO(2,3-D) FURAN	UG/KG	<del></del>				<del></del>		3-0 110				300 110		780 NJ
	UG/KG	<u> </u>				380 NJ		670 NJ			<del></del>	<del></del>		700110
BENZO[B]NAPHTHO[2,1-D] THIOPHENE BENZO[JIFLUORANTHENE	UG/KG	<b></b>			<del></del>	200 140		370 (%)			<del> </del>			<del> </del>
7H-BENZ (DE) ANTHRACEN-7-ONE	UG/KG					310 NJ		950 111	910 NJ	<del></del>	1100 NJ	430 NJ		1500 NJ
CARYOPHYLLENE	UG/KG		<b> </b>			עא טונ		920 M	all No	<del> </del>	1100 NJ	430 NU		1500 NJ
									1000 111		<del></del>	<del></del>		
CHRYSENE, 1-METHYL CHRYSENE, 5-METHYL	UG/KG							580 NJ	1200 NJ	<del> </del>				910 NJ
	UG/KG					270 NJ		290 MJ		ļ				910 143
CHRYSENE, 6-METHYL		<b> </b>				270 NJ		<u> </u>				200 111		1400 NJ
CYCLOPENTA(DEF) PHENANTHRENONE	UG/KG							<u> </u>	1400 NJ		940 NJ 690 NJ			770 NJ
5,12-NAPHTHACENEDIONE	UG/KG		<u> </u>			330 NJ 290 NJ		1200 NJ	1400 NJ		3300 NJ			770 NJ
BENZO[E]PYRENE	UG/KG					290 NJ		1200 NO			3300 NJ			
BENZO(B)TRIPHENYLENE												380 NJ		
BENZEJACEPHENANTHRYLENE	UG/KG													<b></b>
1,2:7,8-DIBENZPHENANTHRENE	UG/KG					260 NJ					<b></b>	480 J		<b></b>
[3,4·9,10] DIBENZPYRENE	UG/KG		L			180 NJ					<b>├</b>			
UNKNOWN ALKENE	UG/KG										ļ			
UNKNOWN CARBOXYLIC ACID	UG/KG										720 J			
UNKNOWN PAHS/NO.	UG/KG					2700 J/11		5300 J/7	5400 J/4		9400 J/1	2100 J/6		8200 J/6
HEXADECANOIC ACID	UG/KG						100 NJ					670 NJ		L
9-OCTADECENE,(E)	UG/KG						170 NJ				l			
1-HEPTADECENE	UG/KG						120 NJ							
D:C-FRIEDOOLEANAN-3-ONE	UG/KG						860 NJ							
9 H-FLUOREN-9-ONE	UG/KG							460 NJ						
CHOLESTEROL .	ÚG/KG						I							
NAPHTHALENE, 2-PHENYL	UG/KG							760 NJ						
1-TETRADECENE	UG/KG									140 NJ				
1-OCTADECENE	UG/KG							-		230 NJ				
1-HEPTADECENE	UG/KG									230 NJ				
3,4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG						T i				550 NJ			
1-EICOSANOL	UG/KG													
TESTOSTERONE .	UG/KG													
TRIPHENYLENE,2-METHYL	ÚGÆG										800 NJ	370 NJ		
VALENCENE	UG/KG													
PESTICIDES POSITION TO THE POSITION TO THE PERSON TO THE P	STEP SHARE SAFE	arabitele dituation	Distriction in a	and the latter of the second	the a silver	a Marke Market of	. The second second	113, 10,445.5	すが、10年代後年代	1. 22. combección	Comment of the last	وعود المال ويوارث ليما والعوارات	arm in armin	August Property
ALPHA-BHC	UG/KG	2 U				- 1	1	-	- 1	-				_
ENDRIN	UG/KG	3.8 U				-	-	_	-	_				182 1 42 J
ENDRIN ALDEHYDE	UG/KG	3.6 U	-			18:31:31			7.17.900 July				<del></del>	56-1502 <b>49.J</b> .50
NDRIN KETONE	UG/KG	3.8 U	-			_			-					150
4,4'-ODT	UG/KG	3.8 U	_						81418.50% J.C.			77.94 to <b>30 d</b> 250		36 U
4'-DDE	UG/KG	38 U		[	- I		- 1	- 1		-		_		

- GS Gudi States Creosoting
  SS surface soil sample
  UGNG micrograms per keogram
  U value is below the reporting limit
  J estimated value
- data is unusable
- · Fresumptive evidence of presence of materix

## TABLE 6 (CONTINUED) SUMMARY OF ORGANIC ANALYTICAL RESULTS SURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

		Bee ground	60.335.444.	independent	4.5 - 1.4 m 4.2 - 4.4 m	et established in the said	SELECTION STORY	re in On-alle	Semples (2)	in to describe sources	SA SECTION AND THE PARTY OF THE	Carrie Service 4	deneral della coll	11,000,000,000
	· · · · · ·	DECEMBER	14/20/20/20	3-1-17 2-13 18	32 2 44 × 440		1	1		T	, Kana Carrie	1		T
			GS14SS		GS16SS	GS17S8	GS18SS	GS19S5	GS20SS	GS20SD*	GS21SS	GS22SS	GS23SS	GS24SS
COMPOUND	UNITS	GS01SS		GS15SS										
OLATELES MATTLE TERRITORISM TO THE TERRITORISM TO T						PER GRANILLAND			83 J	I 230 J	270 J	110 J	130 J	220 J
ACETONE	UG/KG	100 J	75 J	100 J	56 J	260 J	130 J	180 J	- 63 3	230 0	2/03		130 3	220 3
METHYL ACETATE	UG/KG	11 U	-	3 J				3 J					SEL 114: 4:40	L
METHYL ETHYL KETONE	UG/KG	11 U	L <u></u>	1232 a 144	1	5: 28 ca Jaka	1:014 % J 65	247 3.3.3.	<u> </u>	230.5	Sarane G C J. Car	<u> </u>	Ø 0~ (1 € · A+) ?	<del></del>
MISCELLANEOUS VOLATILE ORGANICS								garagie parties,	A. 18. 18. 18. 18. 18. 18. 18. 18. 18. 18					
UNKNOWN COMPOUNDS/NO.	UG/KG	19 J/2	7 J/1		16 J/2	12 J/1	8 J/1			9 J/1	12 J/1	6 J	30 J/3	150 J/2
ACETALDEHYDE	UG/KG		12 NJ	24 NJ		38 NJ		18 NJ		42 NJ	47 NJ	l	14 NJ	İ
BUTANAL	UG/KG		1						L	l			L	L
BUTANAL, 3-METHYL-	UG/KG			1						11 NJ		L		I
HEXANAL	UG/KG		I		31 NJ		20 NJ							1
PENTANAL	UG/KG										7 NJ			
PROPANAL, 2-METHYL-	UG/KG		L		I	7_NJ				7 NJ		L		
XTRACTABLES*	· · · · · · · · · · · · · · · · · · ·	Ann. W 2111	. 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	وي استرد خي رهوم	1 50 mm 6 677	13.72 a 200° s 240	e ter grant to a	्रिकृत- एक्- वर्द्धाः	1966		5	or in wait is	Mariani, u	rain an lagar
3-AND/OR 4-)METHYLPHENOL	lugkig	380 U	1.58**(E.86		si_(K)Hib ()*-3		-		e (jaky) ( / N · · · · )		1924-175		-	Agy
CENAPHTHENE	UG/KG	380 U	<del></del>	<del></del>	<del></del>	<del></del>	<del> </del>	_		<del>  </del>		<del></del>	<del></del>	<del> </del>
CENAPHTHYLENE	UG/KG	380 U	<del>-</del> -	250 J	87 J	270 J				<del>                                     </del>				150 J
NTHRACENE	UG/KG	380 U	-1100 oc	370 J	. B/ J	270 J				<del>  </del>		_	96.7	330 J
BENZALDEHYDE	UG/KG	380 U	AHOO AS	3/0 3		10 304 31.2M			<del></del>	<del>                                     </del>			<del></del>	3303
			2300 (22)		200 1	1000	~			44 J		-	70.7430 J 100	
ENZO(A)ANTHRACENE	UG/KG	380 U	4200	< 1500 ···	260 J			<del>-</del>	58 J	140 J	<del></del>		**** ****	
ENZO(B)FLUORANTHENE 6200/2/00						12900 J.2%			160 J				400	⊕ 2700°3 ≥
ENZO(GHI)PERYLENE	UG/KG	380 U	820	360 J	92 J	290 J	_			-			180 J	330 J
SENZO(K)FLUORANTHENE	UGAKG	380 U	2400	N1100 4		#2700 J	-		170 J	150 J			620	. 2000 J
BENZO-A-PYRENE 621/2100	UG/KG	380 U	2300 × %	413- <b>650</b> Fast	230 J	±1000 ⊊∜;;≥.		-	46 J	47 J		-	380 J	×870.6%.
,1-BIPHENYL	UG/KG	380 U					-			-				
CARBAZOLE	UG/KG	380 U	280 J	250 J	62 J	130 J						-		
CHRYSENE	UG/KG	380 U	3100	1700	4.390 J 21	6:1 <b>600</b> cdefa			98 J	69 J		-	1 1610 FB	in.1100 4505
DIBENZO(A,H)ANTHRACENE 620/210		380 UJ	580 ·	200 J	95 J	330 J		-		<u> </u>			120 J	260 J
DIBENZOFURAN	UG/KG	380 U		-	-		-	-	-					
LUORANTHENE	UG/KG	380 U	· 3100 · · /	93100 as	HU78018F-1	© <b>2600</b> %ind≥t	-		73 J	. 81 J		ائہ	or 520 D. 4	1300 / 5
LUORENE	UG/KG	380 U i				-	_	•	-					-
NDENO(1,2,3-CD)PYRENE	UG/KG	380 UJ	1700	740 Jul	270 J	890	-		63 )	_ 51 J	_		340 J	#3550 J⇒r
-METHYLNAPHTHALENE	UG/KG	380 U	-		-	-	=		-		_			'
APHTHALENE	UG/KG	380 U	-		-	-	<u> </u>			-		-		
HÉNANTHRENE	UG/KG	380 U	320 J	a.1100 /***	310 J	250 J			-	-		_	59 J	
ENTACHLOROPHENOL	UG/KG	970 W			-	-		-				"	-	-
YRENE	UG/KG	380 U	3300 🖘	2200.	J. 540 Jr	1/1700 x 2/2			100 J	63 J			⊅: <b>450</b> ≈ 23	43 <b>1800</b> 445
ISCELLANEOUS EXTRACTABLES	હારમાં ત્યાં ઉપયુક્તિ કહે છે.	Salas in the second	vá děn rokustí	5184 674 G	Out of the last	A POST OF PRO	and the American	bien inga i Feri	Cir scott des	ing the of the charles	9902400 A-230	Mark Control	341 FEV. 3 (F. 167)	4076-0218-80
INKNOWN COMPOUNDS/NO.	UG/KG		5500 J/8 I	12000 J/19	24000 J/17	5900 J/16	3100 J/17	12000 J/25 I	8800 J/17	9100 J/17 I	3100 J/14	1300 J/4		
TIGMAST-4-EN-3-ONE	UG/KG									320 NJ			170 NJ	2600 NJ
NKOWN ALCOHOL	UG/KG							<del></del> -					-,,,,,,,,,	1000.40
NKNOWN KETONES/NO.	UG/KG													<del></del>
ETHYLANTHRACENE	ÚG/KG					<del></del>				<del></del>				
10-ANTHRACENEDIONE	UG/KG			700 NJ	160 J	<del></del>		<del></del>		<del></del>				
HENANTHRENE, 1-METHYL	- 100,10		<del></del>							<del></del>				
HENANTHRENE, 2-METHYL						· · · · · · · · · · · · · · · · · · ·		<del></del>		<del>                                     </del>				
HENANTHRENE, 2,3-DIMETHYL	UG/KG	<b></b>	<del></del>							<del></del>		<del></del>		
HENANTHRENE, 4,5-DIMETHYL	. Journa				<del></del>	·								
NTHRACENE, 2-METHYL								<del></del>						
ENZOFLUORENE	- Heave	<b>_</b>												
	UG/KG	<b>_</b>	4700 4::											
H-BENZO(A)FLUORENE	UG/KG		1700 NJ											
H-BENZO[B]FLUORENE	- 1		I	350 NJ		510 NJ				i				

## TABLE 6 (CONTINUED) SUMMARY OF ORGANIC ANALYTICAL RESULTS SURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

						******		230 00 A= -=	. Camples	2000 1 1 10				1. NO. 1
						公司 医脱氧化物								
COMPOUND	UNITS	GS01SS	G\$14SS	G\$15S8	GS1655	GS1788	GS18SS	GS19SS	GS20SS	GS20SD*	GS21SS	GS22S5	QS23SS	GS24SS
2,2-BINAPHTHALENE		L	·	<u></u>			<del></del>	<del> </del>	<b>↓</b>	<del></del>				<u> </u>
1,2:3,4-DIBENZPYRENE				190 NJ	L	290 NJ		1	.1		Ī			
1,2 4,5-DIBENZPYRENE			L		L			<u> </u>				L		
1-DOCOSENE		L	L							220 NJ				
1-EICOSANOL					l		I	1			230 NJ			
PENTADECANOIC ACID					l	l	1		<u> </u>					L
PERYLENE			710 NJ		200 J	i	i		<u> </u>					260
PYRENE, 1-METHYL			1300 NJ		2000 J	l						l		160 NJ
PYRENE, 2-METHYL	UG/KG		720 NJ		1400 J	L							130 NJ	
PYRENE, 1,3-DIMETHYL	UG/KG					I		Γ	1					
PYRENE, 4-METHYL						1		I	I					
BENZIAJANTHRACENE, 7-METHYL-								I	I				240 NJ	
BENZ(A)ANTHRACENE-7, 12-DIONE				260 NJ			1							
BENZO(B)NAPHTHO(2,3-D) FURAN					· · · · · ·	1			1					
BENZO(B)NAPHTHO(2,1-D) THIOPHENE	UG/KG		550 NJ		1400 J	310 NJ		I						
BENZOUJFLUORANTHENE								· ·					440 NJ	
7H-BENZ [DE] ANTHRACEN-7-ONE	UG/KG		870 NJ	440 NJ	1600 J	390 NJ	1						210 NJ	140 NJ
CARYOPHYLLENE							T	Ì						1600 NJ
CHRYSENE, 1-METHYL				300 NJ				1	t	1				
CHRYSENE, 5-METHYL	<del></del>				2300 J	310 NJ	1			-				
CHRYSENE, 6-METHYL	UG/KG								1	-				
CYCLOPENTA(DEF) PHENANTHRENONE			470 NJ			<u> </u>	†	1					140 NJ	
5,12-NAPHTHACENEDIONE	UG/KG		380 NJ		2600 J	800 NJ		<b></b>	<del>                                     </del>				210 NJ	<u> </u>
BENZO(E)PYRENE	UG/KG		2600 NJ			880 NJ	<del>                                     </del>		<u> </u>				240 NJ	840 NJ
BENZOIBITRIPHENYLENE														
BENZIEJACEPHENANTHRYLENE						f	1	T		-			110 NJ	
1,2:7,8-DIBENZPHENANTHRENE	UG/KG		550 NJ						<b>†</b>	-				
[3,4 9,10] DIBENZPYRENE	UG/KG								<del> </del>	<del>   </del>				
UNKNOWN ALKENE									1	1				200 J
UNKNOWN CARBOXYLIC ACID						t <del></del>			<del>                                     </del>	1				
UNKNOWN PAHS/NO.	UG/KG		5600 J/8	2000 J/4	580 J/1	970 J/3				<del> </del>				1700 J/6
HEXADECANOIC ACID							110 NJ	160 J	130 NU	150 NJ		110 NJ		
9-OCTADECENE,(E)						· · · · · · · · · · · · · · · · · · ·								
1-HEPTADECENE										-				
D:C-FRIEDOOLEANAN-3-ONE					<del></del>		···							
9 H-FLUOREN-9-ONE										<del>                                     </del>		-		
CHOLESTEROL						650 NJ				<del></del>				
NAPHTHALENE, 2-PHENYL	-					210 NJ				-				<del></del>
1-TETRADECENE			- 1			E10 140	-			<del> </del>				
1-OCTADECENE										<del> </del>		-		
1-HEPTADECENE									<del> </del>	<del></del>	-			
3.4-DIHYDROCYCLOPENTA (CD) PYRENE			520 NJ						-	<del> </del>				
1-EICOSANOL			320 140						190 NJ	<del>                                     </del>		<del></del>		
TESTOSTERONE			·	_				390 NJ	1	$\vdash$				
TRIPHENYLENE,2-METHYL			840 NJ				_	330140	<del></del>	<del></del> -				
VALENCENE			040 140		1200 NJ				<del>                                     </del>	<del></del>				
*PECIACIAN					1200 140				1		1			<u> </u>
nestrones and the second	and the second	22 1.0 T. N.	F 54 60 55 000		THE NAME OF THE	1 7 LE 0 R.		h I / . apperta . N s St	deed of old Park	Sa Special Control				
PESTICIDES/PCB	lug/kg I	2 U	Servicia in Principales			CAMPS TOTAL CONTROL	Interior Comme	alleric research class	# 6 8 82 TO 19 12					
		3.8 Ü			-	m v. Brakelinens					**		1.7 NJ	
ENDRIN A DELIVOR	UG/KG					% A B.4 M 1925			<del>-</del>			<u>-</u> .		
ENDRIN ALDEHYDE	UG/KG	3.8 U		27 NJ				-						
ENDRIN KETONE	UG/KG	3.8 U				13c ii 20 J 17c					-		6.9 J	
4,4'-DDT	UG/KG	38 U	10 NJ	30 NJ				-				·-	-	
4,4'-DDE	UG/KG	38U			2.5 J	<del>-</del>		•	4.1 NJ	3.3 NJ				
METHOXYCHLOR	ÜG/KG	20 U	34 NJ		-	-	-	-	-			-	-	-

GS - Guff States Creosoting
SS - surface soil eample
UG/KG - micrograms per kilogram
U - value is below the reporting finit

U - value is below the reporting limit
J - estimated value
R - date is unusable
N - presumptive evidence of presence of material
Sample GS20SD is a duplicate of sample GS20SS shading - elevated concentrations of constituers

# TABLE 7 SUMMARY OF INORGANIC ANALYTICAL RESULTS SUBSURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

		Background	<b>建</b> 数据总统			1.000000000	On-s	ite Samples				zeroza
COMPOUNDS	UNITS	GS01SB	GS02SB	GS03SB	GS04SB	GS05SB	GS06SB	GS07SB	GS08SB	GS09SB	GS10SB	GS11SB
METALS	学に対象	<b>WARRIED</b>		图4.等差别		行為認識的政	and a complete and	<b>。在1960年,</b> 对19				
ALUMINUM	MG/KG	7300 J	7800 J	8600 J	6600 J	5300 J	7700 J	9400 J	8100 J	5800 J	8600 J	9600 J
ANTIMONY	MG/KG	0.59 UJ									0.96 R	1.6 R
ARSENIC	MG/KG	6.7	3.4 J	9.4	1.9	0.91	2.7	2.7	3.5		13	6.6
BARIUM	MG/KG	62	34	63	57	59	48	50	45	29	430	41
BERYLLIUM	MG/KG	0.31	0.36	0.64	0.41	0.27	0.24	0.35	0.59	0.35	1.3	0.49
CADMIUM	MG/KG	0.05 U	i								.XIII 0.87	
CALCIUM	MG/KG	850	170	1400	670	580	990	670	450	540	9400	320
CHROMIUM	MG/KG	12	9.5	14	6.6	4.9	7.5	8	9.2	5.3	19	17
COBALT	MG/KG	3.3	2.6	4.3	1.4	0.78 R	1.2	1.6	9.4	1 R	h / <b>12</b> €	3.2 R
COPPER	MG/KG	6.4	-	·		••		·			40 J	
IRON .	MG/KG	13000	13000	22000 J	7500 J	2300 J	5800 J	8200 J	12000 J	3800	20000	20000
LEAD	MG/KG	8.2 J	6.6	15	9.2	6.7	11	11	. 5.7	8.4	71 图	6.7
MAGNESIUM	MG/KG	580 J	480 J	770	370	220	580	550	370	290 J	2400 J	680 J
MANGANESE	MG/KG	180	200	200	60	16	18	19	1000	9.3	1800	120
NICKEL	MG/KG	5.2 U								<b></b> .	: // (20 %)	
POTASSIUM	MG/KG	320 J	320 J	360	190	170	240	260	350	190 J	.≫0. 970 U	430 J
SELENIUM	MG/KG	0.7	1.4	1.6							1.2	1.1 R
SILVER	MG/KG	0.38	0.38 R	0.49 R			0.28 R		0.32		0.76	1.1 R
SODIUM	MG/KG	290	290	98 J	300	190 J	250	220	46	440	490	350
VANADIUM	MG/KG	22	20	33	14	6.7	17	18	22	8.4	.30	33
ZINC	MG/KG	18	20	23	15	7.3	12	15	25	12	290.	44

### Notes:

GS - Gulf States Creosoting

SB - subsurface soil sample

MG/KG - milligrams per kilogram

U - value is below the reporting limit

J - estimated value

R - data is unusable

shading - elevated concentrations of constituents

# TABLE 7 (CONTINUED) SUMMARY OF INORGANIC ANALYTICAL RESULTS SUBSURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

	Ī											
<del></del>	T	Background			rikupyeye		On	-site Sam	oles		ind, i Turkāsi I	15.00 (10.00 (10.00)) 1
COMPOUNDS	UNITS	GS01SB	GS12SB	GS13SB	GS14SB	GS15SB	GS16SB	GS17SB	GS18SB	GS19SB	GS20SB	GS21SB
METALS	和新的技术	The Edition				(PANTANI)	的复数形式 政治公		Michael Constitution	Printer Beach		
ALUMINUM	MG/KG	7300 J	3700 J	8400 J	4600 J	11000 J	12000 J	6700 J	7700 J	12000 J	4700 J	8600 J
ANTIMONY	MG/KG	0.59 UJ		0.58 R			<u></u>	<u></u>				
ARSENIC	MG/KG	6.7	4.6	4.7	7.4	8.1	7.3	1.5	4.1	8.3	2.3	4.5
BARIUM	MG/KG	62	19	23	54	45	49	61	26	43	24	28
BERYLLIUM	MG/KG	0.31	0.3	0.34	0.48	0.59	0.49	0.39	0.31	0.46	0.3	0.3
CADMIUM	MG/KG	0.05 U	••	-			-		•		1	••
CALCIUM	MG/KG	850	160	400	620	710	390	620	190	190	120	140
CHROMIUM	MG/KG	12	8.5	11	7.1	21	17	7	8.7	12	4.7	14
COBALT	MG/KG	3.3	1.4	1.7	12	2.9	2.1	1.5	3.4	9.6	2.2	1.9
COPPER	MG/KG	6.4				7 J	7.7					3.6
IRON	MG/KG	13000	13000 J	14000	14000	24000	23000	5900	12000	20000	6900	13000
LEAD	MG/KG	8.2 J	6.7	5.1	27	9.2	8.7 J	. 9.1 J	8 J	12 J	6 J	6.8 J
MAGNESIUM	MG/KG	580 J	170	470 J	250 J	640 J	650 J	310 J	470 J	680 J	210 J	420 J
MANGANESE	MG/KG	180	55	82	1000	410	52	57	76	600	140	95
NICKEL	MG/KG											
POTASSIUM	MG/KG	320 J	170	410 J	180 J	490 J	470 J	280 J	220 J	480 J	170 J	300 J
SELENIUM	MG/KG	0.7	1.4			2.3ປ∝	1.7 J			1.7 J		0.95 R
SILVER	MG/KG	0.38	0.47	0.39	0.42	0.65	0.45	0.32 R	0.32 R	0.53	0.25 R	0.4
SODIUM	MG/KG	290	96	310	260	290	290	340	540	280	340	310
VANADIUM	MG/KG	22	26	22	23	36	33	13	18	31	14	21
ZINC	MG/KG	18	8.4	20	20	36	25	12	16	30	11	16

### Notes:

GS - Gulf States Creosoting

SB - subsurface soil sample

MG/KG - milligrams per kilogram

J - value is below the reporting limit

J - estimated value

R - data is unusable

shading - elevated concentrations of constituents

## TABLE 8 SUMMARY OF ORGANIC ANALYTICAL RESULTS SUBSURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

		Background	in application exercisis	density and the	Committee of the second	31.1715.19 · 170.00	entire on On	eite Semples	himbon and the Con-	ian iko didak bibu	Esteropario (Monte	artor o la colonida de colonida.
	<del>†</del>	- Consultation	200 1 4 9 M. MAN 19 1	Lines Francisco	1				7		1	
COMPOUNDS	UNITS	GS01SB	GS02SB	GS03SB	GS04SB		G\$06SB	GS07SB	GS08SB	GS09SB	G\$10SB	GS11SB
VOLATILES ACETONE	IUG/KG		12 J		्रहार <sub>िल्</sub> सिक्ष्यक्रक	10 10 70 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	7,4 ( ) 44 ( ) 45 ( ) 45 ( ) 45 ( ) 45 ( )	 20602423400	99 J	#7181101748686016	78 J	24 J
BENZENE	UG/KG			- 33 5				<del>                                     </del>	***	<del></del>	2 J	
METHYL ETHYL KETONE	UG/KG						<u> </u>		SPECIAL PROPERTY.			
MISCELLANEOUS VOLATILE ORGANICS			7.51 2.5.			VIII I VECE II.	100 S Tart 24 Tart				Section 2020 to 15	*
UNKNOWN COMPOUNDS/NO.	lug/kg		at garaga an m	1	1		8 J/1	. gasangraji da	8 J/1		T 25 J/2	1999.1
EXTRACTABLES (1996)			AND THE WAY	304,281 33875.277241	Net's of such	age artiferation (	189-300-798-15	T 5 5 50 4.	รายปรุงสามารถสา		La North Thirty	55, 50 Mgs.;;;
ANTHRACENE	JUG/KG	400 U	-						74 J		12000	49 J
ACENAPHTHENE		400 U									210 J	
ACENAPHTHYLENE		400 U	•	-	-						6100	
BENZO(A)ANTHRACENE	UG/KG	400 U			<u> </u>			-	120 J	<u></u>	23000	<u></u>
BENZO(B)FLUORANTHENE	UG/KG	400 U		<u> </u>	-		-	47 J	∴ 1000 J →		37000	<u></u>
BENZO(G,H,I)PERYLENE	<del> </del>	400 U			<del></del>	<del></del>	<u> </u>		320 J		5200	
BENZO(K)FLUORANTHENE	UG/KG	400 U				<del>-</del>			∠ to ± 980 J.		26000	48 J
BENZO-A-PYRENE	UG/KG	400 U	L	<u> </u>		- 450 AJ	<del></del>		240 J		25000	
BIS(2-ETHYLHEXYL)PHTHALATE CARBAZOLE	+	400 U	<del></del>		<del></del>	********************************	<del>-</del>	<del></del>	<del></del> -		4.1800 and	<del></del>
CHRYSENE	ÜG/KG	400 U	- <u>-</u>	-:-	-		<del> </del>	<del>"</del>	220 J		35000	<del></del>
DIBENZO(A,H)ANTHRACENE	JUG/KG	400 U	<del></del>	<del></del>			<del></del>	<del>-                                    </del>	150 J		3300	
DIBENZOFURAN	<del> </del>	400 U	<del>-</del>		<del></del>	<del></del>				<del></del>	340 J	<del></del>
FLUORANTHENE	UG/KG	400 U			<del> </del>				200 J		28000	<del></del>
INDENO (1,2,3-CD) PYRENE	UG/KG	400 U							*** 420 CE		20000	
PHENANTHRENE	UG/KG	400 U									2700	
PENTACHLOROPHENOL	1 June	1000 Ü									680 J	
PYRENE	UG/KG	400 U							230 J		37000	<del></del>
MISCELLANEOUS EXTRACTABLES . ACC.	· 建筑物的建筑。46条件,06条件	ng riting galakar.	er Herricania	Laxabeta Maria	Property April 10	ALLESS MARKET OF	State of the second	E4000 (4177)	e nead-Act and area	TO GROWING THE	1.6.25 571 612	1,000,000,000
1-PHENANTHRENECARBOXYLIC ACID	1	92 NJ										
UNKNOWN CARBOXYLIC ACID												86 J
UNKNOWN COMPOUNDS/NO.		840 J/5	220 J/2	250 J/2	<u> </u>			190 J/1	1500 J/7	370 J/3	17000 J/4	530 J/3
UNKNOWN AMIDE	<u> </u>								110 J			
UNKNOWN PAHS/NO.	<u> </u>								560 J/3		65000 J/13	
UNKNOWN PHTHALATES	<del></del>										<b></b>	
PERYLENE	·	اـــــــــــــــــــــــــــــــــــــ							490 NJ		16000 NJ	
PYRENE, 1-METHYL	<del>                                     </del>	<b></b>									9200 NJ	
PYRENE, 5-METHYL		<u> </u>			+						6000 NJ	
9,10-ANTHRACENEDIONE	<del></del>											
BENZ[A]ANTHRACENE, 8-METHYL 7H-BENZ[DE]ANTHRACENEDIONE	<del></del>	L			<b>├</b> ──{						2800 NJ	
7H-BENZIDE ANTHRACEN-7-ONE	<del> </del>										4600 NJ	
11 H-BENZO[B]FLUORENE	<del> </del>										9900 NJ	
BENZO(B)NAPHTHO(2,3-D) THIOPHENE	<del></del>				$\vdash$						4600 NJ	
BENZO[B]NAPHTHO[1,2-D] THIOPHENE	<del> </del>				$\vdash$						4000 140	
BENZOIBITRIPHENYLENE	<del> </del>											
BENZOJETPYRENE	<del>                                     </del>					-					6200 NJ	
1.2-BENZENEDICARBOXYLIC ACID	<del></del>											
2.2'-BINAPHTHALENE											3100 NJ	
CYCLOPENTA (DEF) PHENANTHRENONE	1										4700 NJ	
,2:3,4-DIBENZPYRENE						- · · ·			190 NJ			
,2:7,8-DIBENZPHENANTHRENE											7400 NJ	
8,4-DIHYDROCYCLOPENTA (CD) PYRENE											2800 NJ	
LLIPTICINE												
NAPTHO [1,2,3,4-DEF] CHRYSENE									120 NJ			
,12-NAPHTACENEDIONE			1								4600 NJ	
/ALENCENE												
ESTICIDES/PCB			ALL BURE OF THE	WITH-OWEN.		12 H. 4-14 1/8"		1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	的。 第二章	RAME - 1860	. 100 p. 100 p. 100	
,4-DDT	UG/KG	4 Ü		<del>-</del>							110 N	
,4-DDE	UG/KG	4 U					_=					<del></del>
NDRIN KETONE	UG/KG	4 U									180 N	
SAMMA-BHC (LINDANE)	UG/KG	2.1 U									·	
PC8-1254 (AROCLOR 1254)	UG/KG	40 U	<del></del>		_:_			]				_=_
PC8-1260 (AROCLOR 1260)	UG/KG	40 U			1						••	

- GS Gulf States Creosoling SB subsurface soil sample
- UG/KG micrograms per kilogram
- U value is below the reporting limit
- J estimated value

## TABLE 8 (CONTINUED) SUMMARY OF ORGANIC ANALYTICAL RESULTS SUBSURFACE SOIL SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

	<u> </u>	Background	A 244 44 645	estica bellan	richiel de la lace	etikiya dalah pi eti	≪ . Oπ-siα	Samples ::	CAMPAGE PARTY.	en retetta e	Northern Coppers of th	pán nyezékbek
COMPONIE	UNITS	GS01SB	GS12SB	GS133B	GS14SB	GS15SB	GS163B	GS17SB	GS18SB	GS19SB	GS20SB	GS21SB
COMPOUNDS VOLATILES		USUISB	US 123B	7 G3143B	7 22 1 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2 2	GSISSE	031035					
ACETONE	JUG/KG	39 UJ	44		700 <b>55</b> 4-000 J				-		-	
BENZENE	UGAKG	11 U		<del> </del>			<del></del>	-	<del> </del>	<del> - :</del>	<del></del>	
METHYL ETHYL KETONE	UG/KG	11 UJ		<del> </del>	<del>  </del>		<del>  </del>					
WESTIFE ESTRETATIONS	100.00	1, 52				<b></b>						
MISCELLANEOUS VOLATILE ORGANICS	HOW THE SHAPE	eeric yaariinda	الماسية المرافقة المالية	Salat Salat Salat	iffickles item englichtlich	40-15E 184	· California and Street	BOAT HAR	357516374423V	William Televis	The system is	the deposit of the
UNKNOWN COMPOUNDS/NO.	UG/KG			T	· 8 J/1	20 J/2	T					
EXTRACTABLES CONTROL OF THE PROPERTY	A TO THE SAME	The same of the sa	and of the first	A 4975 (40)	Whitely bull high	Santaka da sa	g Part March	" COPPLE A MARIE	\$45,000000000000000000000000000000000000	87,283 <b></b>	ស្ត្រី <u>ប្រស</u> ៀវទេ	فع الأكبية إلى الرائز
ANTHRACENE	UG/KG	400 U	-	T	9- <b>69</b> 0 45-0							
ACENAPHTHENE	UG/KG	400 U	1									
ACENAPHTHYLENE	UG/KG	400 U	•		330 J							
BENZO(A)ANTHRACENE	UG/KG	400 U		56 J	4/1400 (32/58%)	-						
BENZO(B)FLUORANTHENE	UG/KG	400 U		160 J	2.2400							
BENZO(G,H,I)PERYLENE	UG/KG	400 U			19 540 moderate							
BENZO(K)FLUORANTHENE	UG/KG	400 Ú		170 J	5.2000 pages.		-	43 J				
BENZO-A-PYRENE	UG/KG	400 U			3314003×36-13			ļ <u>-:</u>			<u> </u>	
BIS(2-ETHYLHEXYL)PHTHALATE	UG/KG	400 U							<u> </u>			
CARBAZOLE	UG/KG	400 U	,		210 J	-						
CHRYSENE	UG/KG	400 U	-	91 J	2100 8							
DIBENZO(AH)ANTHRACENE	UG/KG	400 U	-		390		-		-			
DIBENZOFURAN	UG/KG	400 U		<del></del>		-	<u> </u>					
FLUORANTHENE	UG/KG	400 U		130 J	1- 2200 - 1 ES	<del></del>	-					
INDENO (1,2,3-CD) PYRENE	UG/KG	400 U		51 J	201100 EC (10)			-				
PHENANTHRENE	UG/KG	400 U			Q 610 0 128 0 1				<u> </u>			
PENTACHLOROPHENOL	UG/KG	1000 U						ļ <u>.                                 </u>		·		
PYRENE	UG/KG	400 U		89 J	2400	<del></del>	L	ــــــــــــــــــــــــــــــــــــــ	L			<del></del>
I-PHENANTHRENECARBOXYLIC ACID UNKNOWN CARBOXYLIC ACID UNKNOWN COMPOUNDS/NO.	UG/KG UG/KG UG/KG			1000 J/2	3900 J/10	310 J	220 J/2		450 J/2	95 J/1	130 J/1	84 J/1
UNKNOWN AMIDE	UG/KG											
UNKNOWN PAHS/NO.	UG/KG				3200 J/7							
UNKNOWN PHTHALATES	UG/KG								230 J/2			
UNKNOWN PHTHALATES PERYLENE	UG/KG UG/KG				420 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL	UG/KG UG/KG UG/KG								230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL	UG/KG UG/KG UG/KG UG/KG				420 NJ 340 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYAENE, 1-METHYL PYBENE, 5-METHYL 9,10-ANTHRACENEDIONE	UG/KG UG/KG UG/KG UG/KG UG/KG				420 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEDIONE BENZ[A]ANTHRACENE, 8-METHYL	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG				420 NJ 340 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL 9,10-ANTHRACENEDIONE BENZIAJANTHRACENEDIONE BENZIAJANTHRACENEDIONE 7H-BENZIDEJANTHRACENEDIONE	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG				420 NJ 340 NJ 320 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL 9,10-ANTHRACENEDIONE BENZIAJANTHRACENE, 8-METHYL 7-H-BENZIDEJANTHRACENE, 10-METHYL 7-H-BENZIDEJANTHRACEN-7-ONE	UG/KG				420 NJ 340 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEDIONE BENZ[A]ANTHRACENE, 8-METHYL 7H-BENZ[DE]ANTHRACENEDIONE 7H-BENZ[DE]ANTHRACENEDIONE 11 H-BENZOBJELUORENE	UG/KG				420 NJ 340 NJ 320 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL 9,10-ANTHRACENEDIONE BENZIOLANTHRACENE, 8-METHYL 7H-BENZIOEJANTHRACENEDIONE 7H-BENZIOEJANTHRACENEDIONE 7H-BENZIOEJANTHRACEN-7-ONE 11 H-BENZOBJELUORENE BENZOBJNAPHTHO(2,3-D) THIOPHENE	UG/KG				420 NJ 340 NJ 320 NJ 500 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEDIONE BENZ[A]ANTHRACENE, 8-METHYL 7-H-BENZ[DE]ANTHRACENE, 10-NE 11 H-BENZOB[ANTHRACENE, 10-NE 11 H-BENZOB[SHITHRACEN-7-ONE 11 H-BENZOB[SHITHRACEN, 10-NE BENZOB[SNAPHTHO[1,2-D] THIOPHENE BENZOB[NAPHTHO[1,2-D] THIOPHENE	UG/KG				420 NJ 340 NJ 320 NJ 500 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYPRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEJONE BENZ[A]ANTHRACENE, 8-METHYL 7H-BENZIDEJANTHRACENEJOINE 7H-BENZIDEJANTHRACENEJOINE 7H-BENZIDEJANTHRACEN-7-ONE 11 H-BENZO[B]FLUORENE BENZO[B]NAPHTHO[2,3-D] THIOPHENE BENZO[B]NAPHTHO[1,2-D] THIOPHENE BENZO[B]NAPHTHO[1,2-D] THIOPHENE BENZO[B]NAPHTHO[1,2-D] THIOPHENE BENZO[B]NAPHTHO[1,2-D] THIOPHENE	UG/KG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL BENZIAMTHRACENEDIONE BENZIAMTHRACENEDIONE 7H-BENZIDEJAMTHRACENEDIONE 7H-BENZIDEJAMTHRACEN-7-ONE 11 H-BENZOBJETUORENE BENZOBJAMPHTHO[2,2-D] THIOPHENE BENZOBJAMPHTHO[1,2-D] THIOPHENE BENZOBJETURPHENYLENE BENZOBJETURPHENYLENE	UG/KG				420 NJ 340 NJ 320 NJ 500 NJ							
UNKNOWN PHTHALATES PFRYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9 (10-METHYL 9 (1	UG/KG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ				230 J/2			
UNKNOWN PHTHALATES PERYLENE PYRENE, -METHYL PYRENE, -METHYL PYRENE, S-METHYL PYRENE, S-METHYL BENZIAMTHRACENE B-METHYL 7+HERIZIDE JANTHRACENE DIONE 7+HERIZIDE JANTHRACENE DIONE 11 H-BENZOBJANTHRACEN-7-ONE 11 H-BENZOBJANTHRACEN-7-ONE 11 H-BENZOBJANTHROLEN-7-ONE BENZOBJANPHTHO(1,2-O) THIOPHENE BENZOBJANPHTHO(1,2-O) THIOPHENE BENZOBJANPHTHO(1,2-O) THIOPHENE BENZOBJENTEPHENYLENE BENZOBJENTEPHENYLENE 1,2-BENZENEDICARBOXYLIC ACID 2,2-BINAPHTHALENE	UGAKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1800 NJ							
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEDIONE BENZIAJANTHRACENEDIONE BENZIAJANTHRACENEDIONE 7H-BENZIDEJANTHRACENEDIONE 11 H-BENZOBJANTHRACEN-7-ONE 11 H-BENZOBJANTHRACEN-7-ONE BENZOBJENAPHTHO[1,2-D] THIOPHENE BENZOBJENAPHTHALENE OVCLOPENTA (DEF) PHENANTHRENONE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ							
UNKNOWN PHTHALATES PFRYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHRAGENEDIONE BENZ/AJANTHRAGENE, 8-METHYL 7-H-BENZ/DEJANTHRAGENEDIONE 7-H-BENZ/DEJANTHRAGENEDIONE 11 H-BENZO(B)FLUORENE BENZO(B)RAPHTHO(2,3-D) THIOPHENE BENZO(B)RAPHTHO(2,3-D) THIOPHENE BENZO(B)RAPHTHO(1,2-D) THIOPHENE BENZO(B)REDEDICARBOXYLIC ACID 2,2-BINAPHTHALENE CYCLOPENTA (DEF) PHENANTHRENDNE 1,2-3,4-DIBENZPYRENE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1800 NJ							
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEDIONE BENZIAMTHRACENEDIONE BENZIAMTHRACENE, 8-METHYL 7H-BENZIDEJANTHRACEN-7-ONE 11 H-BENZOBJANTHRACEN-7-ONE 11 H-BENZOBJANTHRACEN-7-ONE BENZOBJNAPHTHO[2,2-D] THIOPHENE BENZOBJNAPHTHO[1,2-D] THIOPHENE BENZOBJNAPHTHALENE CYCLOPENTA (DEF) PHENANTHRENDNE 1,2-3,4-DIBENZPHENANTHRENE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 500 NJ 270 NJ 340 NJ 1900 NJ							
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEDIONE BENZIAJANTHRACENEDIONE BENZIAJANTHRACENEDIONE 7H-BENZIOEJANTHRACENEDIONE 11 H-BENZOEJANTHRACEN-7-ONE 11 H-BENZOEJANTHRACEN-7-ONE ENZOEJBINAPHTHO[2,3-D] THIOPHENE BENZOEJBINAPHTHO[1,2-D] THIOPHENE BENZOEJBINAPHTHO[1,2-D] THIOPHENE BENZOEJBINAPHTHO[1,2-D] THIOPHENE BENZOEJBINAPHTHOLE,2-D] THIOPHENE 1,2-BENZENEDICARBOXYLIC ACID 2,2-BINAPHTHALENE CYCLOPENTA (DET) PHENANTHRENDNE 1,2-3-DIBENZPYRENE 1,2-3-DIBENZPHENANTHRENE 2,2-7-B-DIBENZPHENANTHRENE 3,2-7-B-DIBENZPHENANTHRENE 3,2-7-B-DIBENZPHENANTHRENE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ							
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL PHENZOE, DAITHRACENE, 6-METHYL PHENZOE, DAITHRACENEDIONE PH-BENZOE, DAITHRACENEDIONE PHENZOE, DAITHRACENE, 7-ONE 11 H-BENZOE, BANTHRACEN-7-ONE 11 H-BENZOE, BHANTHRACEN-7-ONE 11 H-BENZOE, BHANTHRACEN-7-ONE 11 H-BENZOE, BHANTHRACEN-7-ONE BENZOE, DE PHENZOE, BENZOE, BHANTHRACENE BENZOE, BHANTHRACENE BENZOE, PYRENE 1, 2-8-BINAPHTHALENE CYCLOPENTA (DEP. PHENANTHRENDNE 1, 2-3-4-DIBENZYPHENE 1, 2-1-DIBENZYPHENE	UGAKG				420 NJ 340 NJ 320 NJ 500 NJ 500 NJ 270 NJ 340 NJ 1900 NJ							
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL BENZIANTHRACENEDIONE BENZIANTHRACENEDIONE TH-BENZIDEJANTHRACENE-00NE TH-BENZIDEJANTHRACEN-7-ONE TH-BENZIDEJANTHRACEN-7-ONE TH-BENZIDEJANTHRACEN-7-ONE BENZOIBJANPHTHO[1,3-D] THIOPHENE BENZOIBJANPHTHO[1,3-D] THIOPHENE BENZOIBJANPHTHO[1,2-D] THIOPHENE BENZOIBJANPHTHO[1,2-D] THIOPHENE BENZOIBJANPHTHO[1,3-D] THIOPHENE BENZOIBJANPHTHO[1,3-D] THIOPHENE BENZOIBJANPHTHO[1,3-D] THIOPHENE BENZOIBJANPHTHO[1,2-D] THIOPHENE DENZOIBJANPHTHALENE DENZOIBJANPHTHALENE DYCLOPENTA (DET) PHENANTHRENDNE 1,2-3,4-DIBENZPHENANTHRENE 3,4-DIHYDROCYCLOPENTA (CD) PYRENE ELLITICINE NAPTHO[1,2,3,4-DEF] CHRYSENE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1800 NJ 220 NJ 270 NJ 280 NJ 270 NJ							
UNKNOWN PHTHALATES PFRYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9 (10-METHYL 9 (1	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ						120 M	
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHACENEDIONE BENZIAMTHRACENEDIONE BENZIAMTHRACENE, 8-METHYL 7H-BENZIDEJANTHRACEN-7-ONE 11 H-BENZOBJANTHRACEN-7-ONE 11 H-BENZOBJANTHRACEN-7-ONE BENZOBJANPHTHO[1,2-D] THIOPHENE BENZOBJANPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHENUENE EENZOLEJPYRENE 1,2-BENZENEDICARBOXYLIC ACID 2,2-BINAPHTHALENE CYCLOPENTA (DEF) PHENANTHRENDNE 1,2-3,4-DIBENZPHENANTHRENE 3,4-DIHYDROCYCLOPENTA (CD) PYRENE ELIPTICINE NAPTHO [1,2,3,4-DEF] CHRYSENE 5,12-NAPHTACENEDIONE VALENCENE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ 220 NJ 270 NJ 280 NJ 280 NJ				120 NJ		120 NJ	
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-AMTHRACENEDIONE BENZIAJANTHRACENEDIONE BENZIAJANTHRACENEDIONE 7H-BENZIOEJANTHRACENEDIONE 7H-BENZIOEJANTHRACEN-7-ONE 11 H-BENZOBJENTHRACEN-7-ONE 11 H-BENZOBJENTHRACEN-7-ONE BENZOLBJENTHRACEN-7-ONE 12-BENZOLBJENTHROLI, 2-D THIOPHENE BENZOLBJENAPHTHO[1, 2-D THIOPHENE BENZOLBJENAPHTHO[1, 2-D THIOPHENE BENZOLBJENAPHTHOLI, 2-D THIOPHENE BENZOLBJENAPHTHOLI, 2-D THIOPHENE 1,2-BENZENEDICARBOXYLIC ACID 2,2-BINAPHTHALENE CYCLOPENTA (DET) PHENANTHRENDNE 1,2-3,4-DIBENZPHENANTHRENE 3,4-DIHYDROCYCLOPENTA (CO) PYRENE ELLIPTICINE WAPTHO [1,2,3,4-DEF] CHRYSENE 5,12-NAPHTACENEDIONE VALENCENE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG	4.U	第5 - 416		420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ 220 NJ 270 NJ 290 NJ				120 NJ			
UNKNOWN PHTHALATES PPREVENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 6-METHYL PYRENE BENZOLOBIOATHRACEN-7-ONE 11 H-BENZOLBIAPHTHO(1,2-0) THIOPHENE BENZOLBIAPHTHO(1,2-0) THIOPHENE BENZOLBIAPHTHO(1,2-0) THIOPHENE BENZOLBIAPHTHOLI,2-0) THIOPHENE BENZOLBIAPHTHALENE 12-BENZENEDICARBOXYLIC ACID 12-BENZENEDICARBOXYLIC ACID 12-S-DIBLENTHALENE CYCLOPENTA (DEP) PHENANTHRENDNE 12-3-4-DIBENZPHENE 12-3-4-DIBENZPHENE 12-1-1-CHENERAL S-DIHYDROCYCLOPENTA (CO) PYRENE ELLIPTICINE NAPTHO (1,2,3-4-DEF) CHRYSENE 5-12-MAPHTACENEDIONE VALENCINE VALENCINE VALENCINE PERITCIDES/PCS	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG	4 U			420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ 220 NJ 270 NJ 280 NJ 280 NJ				120 NJ		230°446,048	
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL J. 1-METHYL J. 1-MENZONEDIONE BENZIAMTHRACENEDIONE BENZIAMTHRACENEDIONE TH-BENZIDEJAMTHRACEN-7-ONE 11 H-BENZOBJELORENE BENZOBJAMPHTHO[2,3-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHENE 1,2-BENZENEDICARBOXYLIC ACID 2,2-BINAPHTHALENE CYCLOPENTA (DET) PHENANTHRENDNE 1,2-3,4-DIBENZPHENANTHRENE 3,4-DIHYDROCYCLOPENTA (CD) PYRENE ELIPTICINE VAPTHO[1,2,3,4-DEF] CHRYSENE 5,12-NAPHTACENEDIONE VALENCENE PEBTICIDES/PCB.	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG				420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ 220 NJ 290 NJ 290 NJ		-	_=	120 NJ		zarwą jesk 	
UNKNOWN PHTHALATES PFRYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL 9, 10-ANTHRACENEDIONE BENZIAJANTHRACENEDIONE BENZIAJANTHRACENEDIONE BENZIAJANTHRACENEDIONE 77-BENZIOEJANTHRACENEDIONE 11 H-BENZOEJANTHRACEN-7-ONE 11 H-BENZOEJBANTHRACEN-7-ONE BENZOEJBINAPHTHO[2,3-D] THIOPHENE BENZOEJBINAPHTHO[1,2-D] THIOPHENE BENZOEJBINAPHTHO[1,2-D] THIOPHENE BENZOEJBINAPHTHOLE,3-D THIOPHENE 1,2-BENZENEDICARBOXYLIC ACID 2,2-BINAPHTHALENE CYCLOPENTA (DEP) PHENANTHRENDNE 1,2-3-DIBENZPHENANTHRENE 1,2-3-DIBENZPHENANTHRENE 1,2-3-DIBENZPHENANTHRENE BLIJPTICINE MAPTHO [1,2,3,4-DEF] CHRYSENE 5,12-NAPHTACENEDIONE 7-AL-DIDENZPHENE 1,4-DDT 1,4-DDT 1,4-DDT 1,4-DDT 1-DDRINKETONE	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG	4 U 4 U 4 U		·, 6.9	420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ 270 NJ 290 NJ 270 NJ 290 NJ		 		120 NJ		2.487424 (3.144) 	
UNKNOWN PHTHALATES PERYLENE PYRENE, 1-METHYL PYRENE, 1-METHYL PYRENE, 5-METHYL PYRENE, 5-METHYL J. 1-METHYL J. 1-MENZONEDIONE BENZIAMTHRACENEDIONE BENZIAMTHRACENEDIONE TH-BENZIDEJAMTHRACEN-7-ONE 11 H-BENZOBJELORENE BENZOBJAMPHTHO[2,3-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHTHO[1,2-D] THIOPHENE BENZOBJENPHENE 1,2-BENZENEDICARBOXYLIC ACID 2,2-BINAPHTHALENE CYCLOPENTA (DET) PHENANTHRENDNE 1,2-3,4-DIBENZPHENANTHRENE 3,4-DIHYDROCYCLOPENTA (CD) PYRENE ELIPTICINE VAPTHO[1,2,3,4-DEF] CHRYSENE 5,12-NAPHTACENEDIONE VALENCENE PEBTICIDES/PCB.	UGKG UGKG UGKG UGKG UGKG UGKG UGKG UGKG	4 U		6.9	420 NJ 340 NJ 320 NJ 500 NJ 270 NJ 340 NJ 1900 NJ 270 NJ 290 NJ 290 NJ		  		120 NJ		######################################	====

GS - Gulf States Crossoling
SB - subsurface soil sample
UG/KG - micrograms per kilogram
U - value is below the reporting amit

J - estimated value shading - elevated concentrations of constituents

# TABLE 9 SUMMARY OF FIELD PARAMETERS FOR GROUNDWATER SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

Sample Number	pH	Conductivity (µS/cm)	Turbidity (NTU)	Temperature °C
GS01GW	5.41	218	5.97	18.5
GS02GW	4.95	724	3.7	18.4
GS03GW	5.93	180	9.84	17.5
GS04GW	5.81	59	11	18.9
GS05GW	5.32	608	6.23	17.6

### Notes:

GS - Gulf States Creosoting

GW - Groundwater sample

 $(\mu S/cm)$  - microsiemens per centimeter

NTU - Nephelometric turbidity units

°Ć - degrees Celsius

# TABLE 10 SUMMARY OF INORGANIC ANALYTICAL RESULTS GROUNDWATER SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

		Background	1	On-site S	amples	
COMPOUND	UNITS	GS01GW	GS02GW	GS03GW	GS04GW	GS05GW
METALS					<b>建筑建筑</b>	912 A.C.
ALUMINUM	UG/L	320 UJ	450 ປ	850 J.	1700 J	370 J
BARIUM	UG/L	20	120	29	26	32
BERYLLIUM	UG/L	0.11	0.14	0.11		0.9
CALCIUM	UG/L	4600	16000	3600	3500	15000
CHROMIUM	_UG/L	1.8 R	0.92	1.8	1.6	
COBALT	UG/L	1.2	1.2 R			10
IRON	UG/L	1500	1400	2100	1900	12000
MAGNESIUM	UG/L	1500	5500	940	930	***- 8200.::
MANGANESE	UG/L	23		1.10	30	- 土。320小
POTASSIUM	UG/L	520	1500	870	≦ <del>7</del> 21800 : □	1900
SODIUM	UG/L	40000	120000	29000	5000	86000
VANADIUM	UG/L	0.6 U	0.82 R	<b>是35%2</b> 4.	1.9	
ZINC	UG/L	11 U	40,00261367	<b>₩.::</b> -::18		42::

## Notes:

GS - Gulf States Creosoting

GW - groundwater sample

UG/L - micrograms per liter

U - value is below the reporting limit

- estimated value

R - data is unusable

shading - elevated concentrations of constituents

# TABLE 11 SUMMARY OF ORGANIC ANALYTICAL RESULTS GROUNDWATER SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

	-Background	编数别介标的	On-site	Samples 🛸		
UNITS	GS01GW	GS02GW	GS03GW	GS04GW	GS05GW	
		是是建筑的		第6亿日接销售信息		
UG/L	6.1 UJ			7.2 J		
					-	
		造物作品(学		angraf elasiyi		
UG/L			1.3 NJ			
			adalisa see			
Filosofia de Companya de C Transporta de Companya de C	A THE STATE OF THE	er en	aritie Company (1900)	<u> 1984-998 - 1984-988 - 19</u>		
UG/L	2.2 NJ					
UG/L		2.2 NJ	2.1 NJ		1	
UG/L					4.7 J/2	
Same and the same a	outra dimensi	12.30(15.37(2))			第15章 第16章 <del>第16章 第</del>	
	UG/L UG/L UG/L UG/L UG/L UG/L	UG/L C.2 NJ UG/L UG/L UG/L UG/L UG/L UG/L	UNITS   GS01GW   GS02GW   UG/L   6.1   UJ	UNITS   GS01GW   GS02GW   GS03GW   UG/L   6.1   UJ	US/L   C.2 NJ   C.3 NJ   US/L   C.4 NJ   US/L   US/L   C.4 NJ   US/L   C.4 NJ   US/L   US/L	

### Notes:

GS - Gulf States Creosoting

GW - groundwater sample

UG/L - micrograms per liter

U - value is below the reporting limit

J - estimated value

N - presumed presence of constituent

shading - elevated concentrations of constituents

# TABLE 12 SUMMARY OF INORGANIC ANALYTICAL RESULTS SEDIMENT SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

		POWER STATES		Marsh Sample:	488435-EM64		Pearl	River	Prairie Branch
			Downstream Samples				S Control	Control	
COMPOUND	UNITS	G\$01SD	GS02SD	GS03SD	GS04SD	GS05SD	GS06SD	GS07SD	GS08SD
METALS		# <b>#</b> ##################################	SENERAL SERVICES	<b>公会被推图</b> 其实	AND THE STREET		er til flanken	AND THE MESSAGE	etalan etalah etal
ALUMINUM	MG/KG	8200 J	4700 J	16000 J	9100 J	4900 J		140 J	510 J
ANTIMONY	MG/KG	12 UJ				2.1 R	0.7 R		
ARSENIC	MG/KG	5.7	1.8 R	7.9 R	5.5	2.9 R			
BARIUM	MG/KG	89	. 77		120	71	6.8	3.2	. 15
BERYLLIUM	MG/KG	0.39	0.56	23.50 A.3.55	1.2 A	0.51			0.1
CALCIUM	MG/KG	1500	880	2000	2000	1200	250		180
CHROMIUM	MG/KG	13	7.8	24	15	8.4			1.4
COBALT	MG/KG	4.4	4.7	12	11	5.5	0.61 R	0.47	1
IRON	MG/KG	15000	6400	24000	13000	8400	950	340	2200 J
LEAD	MG/KG	12 J	20 J	41 J	21	12	1.8	1.2 J	3.6
MAGNESIUM	MG/KG	1000 J	410 J	1200 J	1000 J	660 J	64 J		81
MANGANESE	MG/KG	340	500	720	280	320	52	53	44
POTASSIUM	MG/KG	700 J	410 J	1100 J	811 J	480 J	90 J	52 J	56
SELENIUM	MG/KG	1.9 R		4.2		•			
SILVER	MG/KG	0.68 R	0.45 R	1.1 R	1.1 R	0.71 R	0.24 R		
SODIUM	MG/KG	690	580	1500	1200	600	330	330	68
VANADIUM	MG/KG	24	13	41	23	· 13	1.6 R	0.57 R	2.4
ZINC	MG/KG	36	33	. 130. ↔	84	39	3.6		4.6
T.O.C.	%	2	2.1	9.5	NR	1.8		NR	<b></b>

### Notes:

SP - Sonford Products

SS - surface soil sample

MG/KG - milligrams per kilogram

U - value is below the reporting limit

J - estimated value

R - data is unusable

NR - no reading

shading - elevated concentrations of constituents

T.O.C. - Total Organic Carbon (%)

## TABLE 13 SUMMARY OF ORGANIC ANALYTICAL RESULTS SEDIMENT SAMPLES GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

			1815 F. V	Harab Quanta	<b>B</b> OOK N 1994	State is a discovery	83 di St. Pearl	Phone	Prairie Branc
					ream Semples			Downstream	Control
and a supplied	Liver		GS02SD	GS03SD	GS04SD	GSO5SD	GS06SD	OS07SD	GS08SD
COMPOUNDS	UNITS	GS01SD							
/OLATILES	lug/kg l	94 J	120		520 J	51 J	<u>ე, ათ წე⊹ე ე-ე.</u> <b>22 J</b>	49 J	29 J
	UG/KG	94 J 5 J	130 J	51 J		- 313	13 U	. 49 J	29 J
CARBON DISULFIDE METHYL ETHYL KETONE	UG/KG	37 UJ	<del></del>	(300) NO (30)	68 J	<del> </del>	13 UJ	<del></del>	14 U
MEINTLEINTLREIONE	UG/KG	37.03	L	<del></del>	1	1	1303	I	14 0
MISCELLANEOUS VOLATILE ORGANICS		31 37873	ija i na kara	प्रशासीकी करते थी।			क्षानुबन्धः विकास सम्बद्धाः । १९४१	West Million o	
JNKNOWN COMOUNDS/NO.	UG/KG			<del> </del>	180 J/1 71 NJ	33 J			8 J
ACETALDFHYDE	UG/KG			٠	/ / Nu	L	l	<u> </u>	
1000 A 40 A 10 A 10 A		1.00	Grand Control of the			and the state of	. 42		A
XTRACTABLES					I=3866 . <b>5</b>		080	TO THE MORNING	0.8 t
CENAPHTHENE	UG/KG	0.5 J	1 J				*** *		
ACENAPHTYLYENE	UG/KG	3	2_		25 6 8 8 8 4 8 5 4 1.	7 7	0.8 U		0.8 U
ANTHRACENE	UG/KG	3			75,41 2/52 <b>34</b> ,516.5		2 U		0.8 U
BENZALDEHYDE	UG/KG	89 J		110 J	130 J		330 U		330 U
BENZO(A)ANTHRACENE	UG/KG	1 J	14 (3,00 (3995) <b>7</b> (3,0			12 🤄	5 U	<del></del>	330 U
BENZO(B)FLUORANTHENE	UG/KG	92 J	<del></del>	> 540 J	<del> </del>	-	330 U	<del>-</del>	
BENZO(GHI)PERYLENE	UG/KG UG/KG	92 J 440 U	<del></del>	√ 560 J		<del>-</del>	330 U 330 U		330 U
BENZO(K)FLUORANTHENE									330 U
BENZO-A-PYRENE	UG/KG	2	3 2 1 10 10 1	120	17	16	4 U 7 U		0.8 U
CHRYSENE	UG/KG	1 J				26			2 U
DIBENZO(A.H)ANTHRACENE	UG/KG	2	1 J		50.	34	1 12 U		0.8 U
LUORANTHENE LUORENE	UG/KG	0.5 J	and the sh≥test shortest wh2 a c		: : : : : : : : : : : : : : : : : : :		0.3 J		
		440 U		180 J	PASECALATINEDO, 11 OC.	1 J			0.8 U
NDENO (1,2,3-CD) PYRENE	UG/KG UG/KG	0.6 J					330 U 0.4 J		330 U
2-METHYLNAPHTHALENE	UG/KG	. 0.6 J	0.9 J		2 J	0.8 J		0.4 J	0.8 U
PENTACHLOROPHENOL	UG/KG	24 U	0.8 3	10.00 Sec. 314.00	2013 254 B B	0.9 3	0.5 J 13 U	0.4 J	0.4 J 13 U
PHENANTHRENE	UG/KG	2 J	জনকার কর্মকার জনকার কর্মকার		4 * 872 G. <b>11</b> ; 15c	<del></del>	4	<del></del>	90
PYRENE	UG/KG	2 2			33 - 33 - 3		10 U		2 U
Thene	JUGANG		A- 10010-10-	ga v ver <b>seu</b> , ev	1.5 - 2 - 3 - 3 - 3 - 3	1 7	100		20
MISCELLANEOUS EXTRACTABLES :	and the second second	day of contract Street	at the same of the same of	Access to the second	. A S 25 26 26 2 . A S.	mar and a second	C. C. S. D. V. (2011) 1848 64	NAME OF STREET	a Reidiler
IEXADECANOIC ACID	lug/kg I	420 NJ	ility A. C. (C. 1811 V. S. 1811)	1	ALC PLANTAGE PROPERTY OF	35 / Aug 1 / A	Salar Salahan Salar Salar		
-HEXADECENE	UG/KG	540 NJ		1000 NJ				<b></b>	
RGOST -5-N-3-OL	UG/KG	1500 NJ		1000 140	1700 NJ		<u> </u>		
GAMMASITOSTEROL	UG/KG	1400 NJ	1400 NJ		1700 150				
TIGMASTEROL	UG/KG	1100 NJ	1400 140						
TIGMAST-4-EN-3-ONE	UG/KG	730 NJ		1400 NJ	2200 NJ	BOO NJ			
INKNOWN COMPOUNDS/NO.	UG/KG	23000 J/23		54000 J	54000 J/24	140000 J/25	61 J/1	69 J/1	170 J/3
INIDENTIFIED COMPOUNDS/NO.	UG/KG	92000 J/28	92000 J	34000 3	34000 3/24	140000 3/25	01.0/1	09 1/1	1/0 3/.
ENZO[E]PYRENE	UG/KG	32000 3120	92000 J	590 NJ					
CHOLESTANOL	UG/KG			350 NJ	1600 NJ				
INKNOWN ALCOHOL	UG/KG				9200 J				
NKNOWN ALCOHOL	UG/KG		ļ	<b></b>	9200 J 1800 J				· · · · · · · · · · · · · · · · · · ·
ENZENE, 1-METHYL-2-ISOPROPYL	UG/KG			ļ	1000 3	1200 NJ			
NKNOWN CARBOXYLIC ACID	UG/KG			<del></del>		500 J			
HENANTHRENONE DERIVATIVE	UG/KG			<del> </del>		2500 NJ			
ESTICIDE8/PCB	IUG/KG I	7.3 U	7.0	K07016, 54.	CHANGE OF BUILDING	1878	angis masukita. 4.1 U		4.11
ETA-BHC	UG/KG	7.3 U	AND 100 100 100 100 100 100 100 100 100 10						4.1 U
	UG/KG	7.3 U		2.5			2.1 U		2.1 U
NDOSULFAN II (BETA)			6.5 J	<del> </del>	<del>-</del>		4.1 U	<u> </u>	4.1 U
NDOSULFAN SULFATE	UG/KG	7.3 U	10 N	14 NJ	-		4.1 U		4.1 U

GS - Gulf States Creosoting

SD - sediment sample |
UG/KG - micrograms per kilegram
U - value is below the reporting limit
J - estimated value

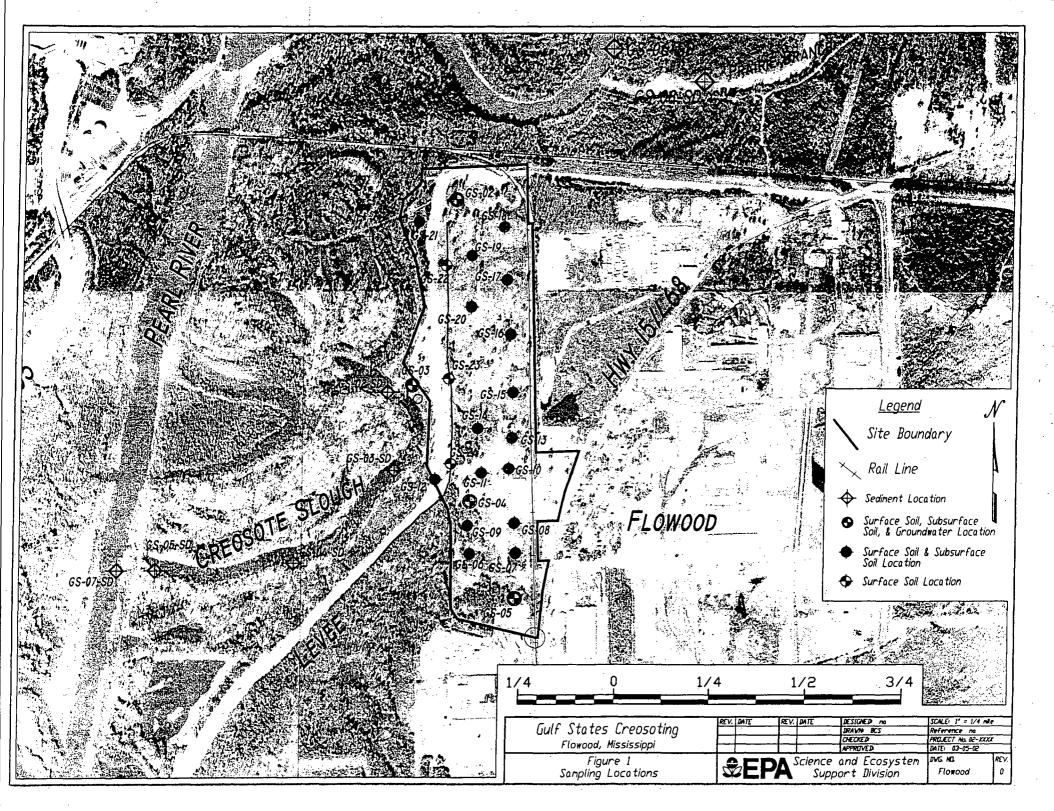
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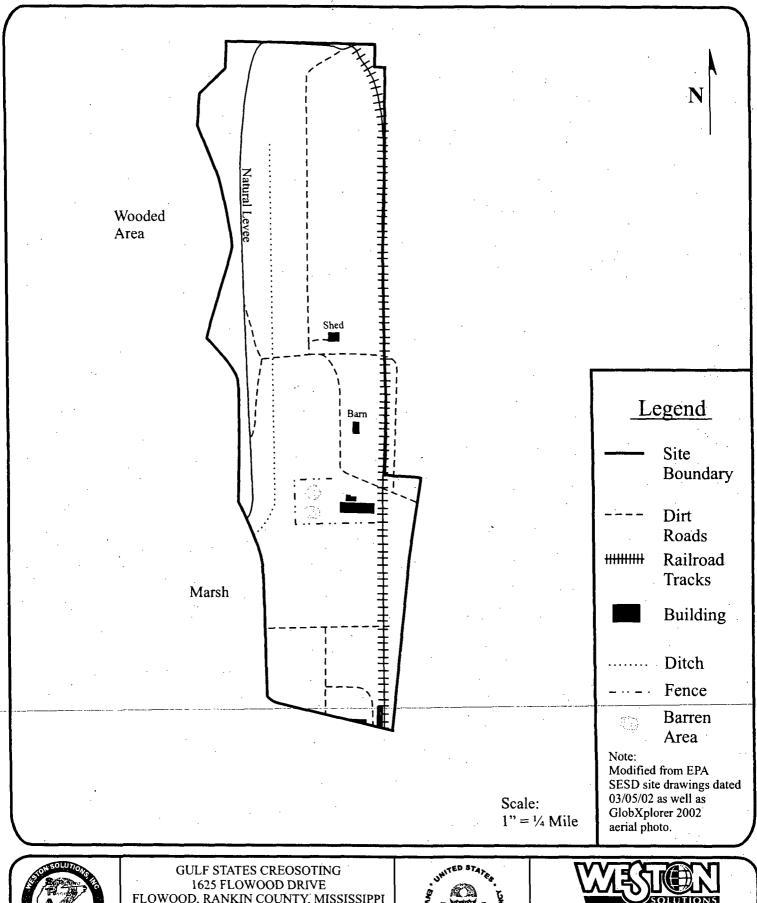
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## APPENDIX B

**FIGURES** 

2 Pages







FLOWOOD, RANKIN COUNTY, MISSISSIPPI

SITE LAYOUT MAP FIGURE 2





12587-001-001-0133-00 4W-02-03-A-003 MSN000407423

## U.S. EPA REGION IV

# SDMS

Unscannable Material Target Sheet

DocID: 10642417	Site ID: <u>msN000407423</u>
Site Name: Duly States	Crossting Campany
Nature of Material:	
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Photos:	CD-ROM:
Blueprints:	Oversized Report:
Slides:	Log Book:
Other (describe): Madeius	map
Amount of material:	
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Reference 3 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423

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## **Coordinate Format**

D/M/S



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NAD83/WGS84



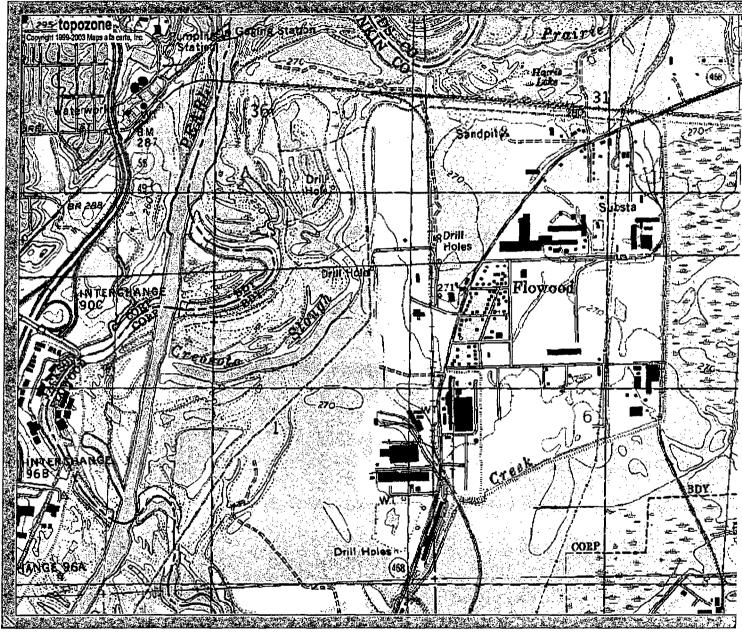
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**USGS Jackson (MS) Topo Map** 

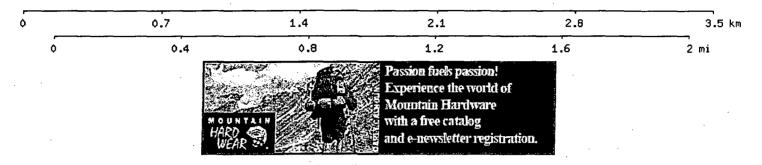
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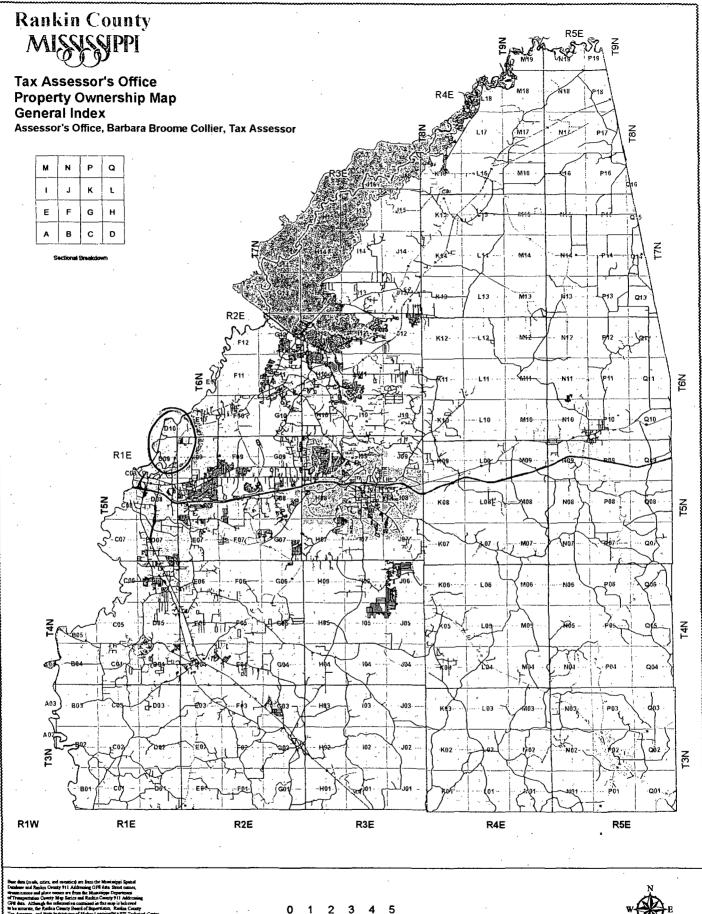
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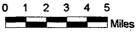
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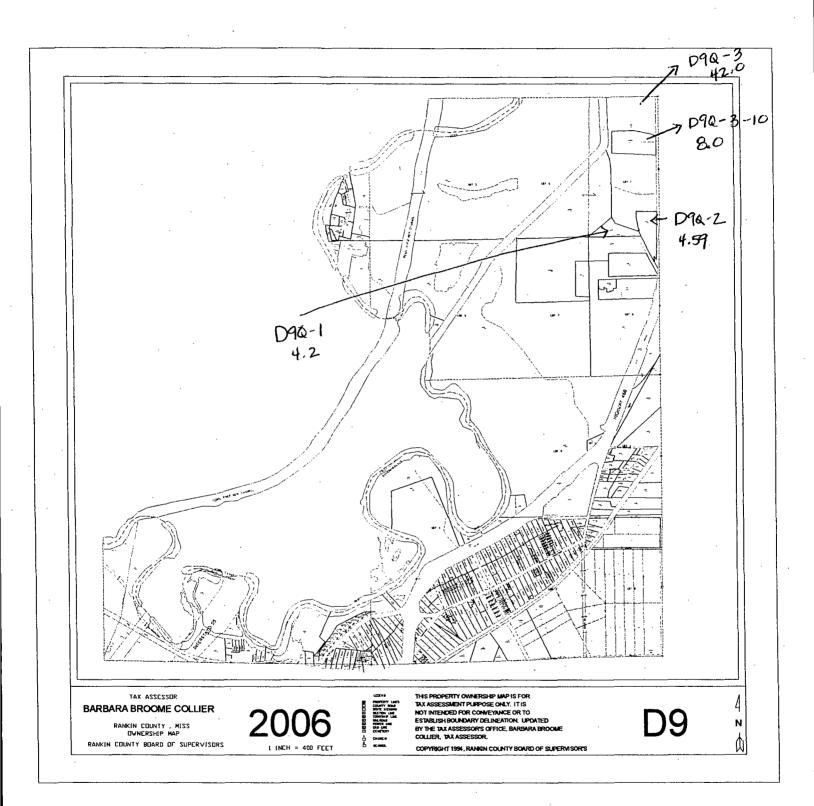
Reference 4
Gulf States Creosoting Company
Flowood, Rankin Co., Mississippi
MSN000407423

Project Note					
Date: June 2, 2008  Gulf State Creosoting Company Flowood, Rankin County, Mississippi TDD Number: TNA-05-003-0045					
Organization: T N & Associates, Inc., Reg. 4 EPA START Contract Name: Stacy Kowalski Signature:	· · · · · · · · · · · · · · · · · · ·				
Subject: Rankin County Tax Assesso	r Information for Gulf States Cre	osoting Company			
The enclosed information contains in Assessor's Office. The subject proper parcel. The information was accessed	erty is the former Gulf States Cre	osoting Company			
	•				
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RE (x) None () Phone	SPONSE REQUIRED call ( ) Memo ( ) Letter	( ) Report			
cc: (x) File ( ) Project Manag	ger ()Principal Investigator	( ) Other (specify)			









PPIN	Parcel Number	/	Section	Township	Range	
049712	D09Q000003 00010	D90-3-10	01	05	. 01	
Assessed Owne			Appraised Va	alues		
CONSTEEL CO IN	iC		Land Value		200,000.00	
P O BOX 6175			Improvement	Value	427,260.00	
JACKSON MS392	88		Total		627,260.00	
Location		<del></del>	Assessed Va	lues		
	DD	<del></del>	Land Value		30,000.00	
1625 FLOWOOD			Improvement	Value	64,089.00	
Legal Descripti			Total		94,089.00	
	' X 416.64 X 432.86 2 IN NE4 NE4(PART		Building Info.			
OF LOT 1)		•	Year Built_	1995		
DB 708 PG 0084	0080294	·	Base Area	20698		
TAXES	Due	Paid	Adjusted Area	3	20760	
County	0.00	0.00	Structure Cod	le	039	
City	0.00	0.00	IDood Info		· · · · · · · · · · · · · · · · · · ·	
School	0.00	0.00	Book			
Total	0.00	0.00	Page			
Paid By		0.00	Date ^	·	0/0/0	
Paid - Receipt #			Miscellaneo	us		
Date Paid	0/0/0		<b>1</b>			
# of Payments	0	<del></del>	1			
Penalty Paid	0	<del></del>	1			

Close

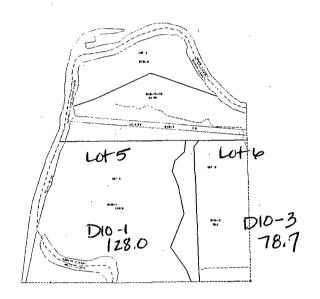
PPIN	Parcel Number		Section	Township	Range		
005816	D09Q000003 00000 / T	90-3	01	05	01		
Assessed Own	ner		Appraised Va	lues			
AVERY LEAD TE	RACK LLC		Land Value		12,900.00		
1 SPRING LAKE	POINTE		Improvement	Value	5,590.00		
PEARL MS3920	8		Total		18,490.00		
Location		<del></del>	Assessed Val	ues			
		<u> </u>	Land Value		1,936.00		
0 FLOWOOD DI			Improvement	Value	839.00		
Legal Descrip			Total		2,775.00		
	IN LOT 1 (E2 NE4) & DO STRIP IN NE COR NE4		Building Info.				
SE4			Year Built		1950		
DB 166 PG 400	0 0060559		Base Area		3740		
TAXES	Due	Paid	Adjusted Area		3754		
County	115.54	115.54	Structure Cod	e .	B42		
City	55.50	55.50	Dood Tota				
School	141.14	141.14	( Rook		731		
Total	312.18	312.18	Page	·	621		
Paid By				Date - 5/15/199			
Paid - Receipt	# Y		Miscellaneou	ıs			
Date Paid	1/17/2008	•	1				
# of Payments	# of Payments 1			····			
Penalty Paid	0		<u> </u>				

Close

6/2/2008

PPIN	Parcel Number		Section	Township	Range	
005815	D09Q000002 00000	DAQ-2)	01	05	01	
Assessed Own	er		Appraised Va	ilues		
CRC PROPERTIE	S LLC		Land Value		114,750.00	
1365 FLOWOOD	DR		Improvement	Value	542,450.00	
FLOWOOD MS3	9232		Total		657,200.00	
l continu	<del></del>	,	Assessed Val	lues		
Location			Land Value		17,213.00	
1353 FLOWOOD		<del></del>	Improvement	Value	81,368.00	
Legal Descript			Total 98,581			
	C IN SE COR LOT 1 E COR LOT 8 (NE SE)		Building Info.			
W/S OF R/R			Year Built 1979			
DB 397 PG 125	0112679	<del></del>	Base Area			
TAXES	Due	Paid	Adjusted Area	·	18144	
County	4,070.41	4,070.41	Structure Cod	<u>e</u>	039	
City	1,971.62	1,971.62	III DAAA TEEA			
School	5,013.83	5,013.83				
Total	11,055.86	11,055.86	Page			
Paid By CRC PROPERTIES LLC			Date 0/0/			
Paid - Receipt # Y			Miscellaneou	us	<del></del>	
Date Paid	1/29/2008	1				
# of Payments 1			<u> </u>		ļ	
Penalty Paid	0		]		<u></u>	

Close



ND22322A XAT

#### BARBARA BROOME COLLIER

RANKIN COUNTY, MISS

RANKIN COUNTY BOARD OF SUPERVISORS



THIS PROPERTY OWNERSHIP MAP IS FOR TAX ASSESSMENT PURPOSE ONLY. IT IS NOT INTENDED FOR CONVEYINGE OR TO ESTBELISH BOUNDARY DELINEATION. UPDATED BY THE TUX ASSESSORS OFFICE, BARBARA BROOME COLLIER, TAX ASSESSOR.

D10

PPIN	Parcel Number		Section	Township	Range		
005820	D10 000003 00000	210-3)	36	06	01		
Assessed Owne	<u> </u>	/	Appraised V	alues			
AVERY LEAD TRA	CK LLC		Land Value		39,550.00		
1 SPRING LAKE P	OINTE		Improvement	: Value	97,840.00		
PEARL MS39208			Total		137,390.00		
Location		· .	Assessed Va	lues			
1649 FLOWOOD	DD.		Land Value		5,933.00		
		<u>.</u>	Improvement	. Value	14,676.00		
Legal Description	5 5&6 SEC 36-5-1 E		Total		20,609.00		
OF LEVEE	5 5&6 SEC 36-5-1 E		Building Info.				
DB 166 PG 400 0			Year Built	1993			
DB 617 PG 0064	0022291		Base Area		6075		
TAXES	Due	Paid	Adjusted Are	a	6075		
County	850.95	850.95	Structure Co	de	039		
City	412.18	412.18	IIDaad Infa				
School	1,048.17	1,048.17	Book		731		
Total	2,311.30	2,311.30	Page	<u> </u>	621		
Paid By AVERY LEAD TRACK LLC			Date 5/15/199				
Paid - Receipt #	<del></del>		Miscellaneo	ous	<del></del>		
Date Paid	1/17/2008		1				
# of Payments 1			<u> </u>				
Penalty Paid	0		]		<u> </u>		

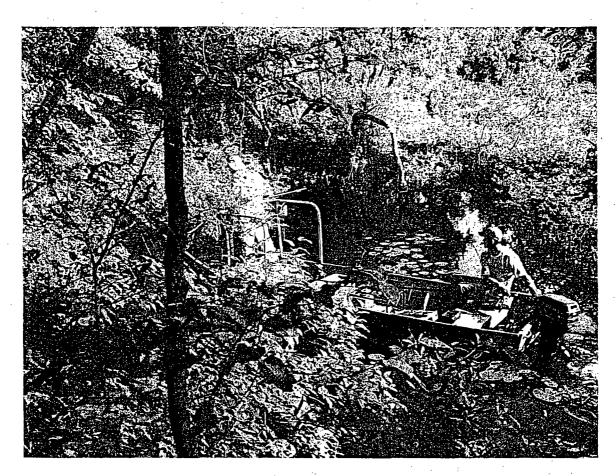
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Reference 5 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423

# **CREOSOTE SLOUGH**

# Water, Sediment and Fish Sampling

Associated with Gulf State Creosoting Co., Inc. Flowood, Mississippi





U.S. EPA, Region 4
Science and Ecosystem Support Division
Ecological Assessment Branch
Athens, Georgia
November 2006

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West 1

#### 1.0 INTRODUCTION

### 1.1 BACKGROUND

The United States Environmental Protection Agency (EPA), Region 4, at the request of the Mississippi Department of Environmental Quality (MDEQ), asked the Science and Ecosystem Support Division (SESD) to assess fish in the area of the former Gulf States Creosoting Company. Since many people consume fish from Creosote Slough, the potential presence of polycyclic aromatic hydrocarbons (PAHs) in fish tissue had become a concern of the MDEQ.

The 141-acre tract, formerly called the Gulf States Creosoting Company, is located at 1625 Flowood Drive (Mississippi Hwy 468), Flowood, Rankin County, Mississippi. The facility lies within an area comprised of mixed industrial, commercial and residential applications (US EPA 2003). The property is currently being used as a horse farm and as a site for ConSteel Co., Inc. Currently, the southernmost building on the property is owned by ConSteel Co., Inc. (Figure 1), which is a steel fabricator and erector that constructs concrete forms, and concrete reinforcement accessories, as well as steel processing and fabrication equipment. The northernmost structures (i.e., metal barn and shed) are used to house tractors and helicopters for a local news station and are part of the horse farm (Figure 2). The horse farm has been operating on the property for approximately 9 years. The facility is bounded by railroad tracks to the north and east, with adjacent businesses to the south, and marshland/tributaries and the Pearl River to the west. The Creosote Slough side of the facility is bounded by a large levee which is fenced and locked (Figure 3). Creosote Slough can be accessed without going through the fenced area by way of the Pearl River.

Gulf States Creosoting Company owned the property as early as 1929 and operated it as a wood treating facility until the mid 1950s. During Gulf States operations, railroad cross ties were treated at the facility with coal-tar creosote and transported on and off-site by means of railroad box cars (US EPA 2003).

Prior to sampling, a site reconnaissance was completed in April, 2006, with assistance from Richard Ball of MDEQ during which time projected sampling stations were chosen. During the reconnaissance, a seep was observed and photographed on the eastern side of the slough near Station CS-04 (Map 1, and Figure 7).

#### 1.2 OBJECTIVE

The primary objective of this survey was to collect a representative set of fish, sediment and water samples within Creosote Slough, which is adjacent to the former Gulf States Creosoting Company, and analyze them for polycyclic aromatic hydrocarbons (PAHs).

#### 2.0 MATERIALS AND METHODS

#### 2.1 SAMPLING LOCATONS

The study area was comprised of a reference station (CS-01) and three other stations (CS-02, CS-03, and CS-04) that covered the area from the railroad tracks on the north end of the property to the terminus of the slough at the Pearl River (Map 1).

#### 2.2 SAMPLING LOGISTICS

The study was initiated and completed during the week of May 29, 2006. To collect the required information, the sampling was conducted when the water-level conditions were low and when there had been no major releases from the Ross Barnett Reservoir to the north of the site.

Personnel from the MDEQ assisted EPA with the collection of water, sediment and fish samples in accordance with the *Ecological Assessment Standard Operating Procedures and Quality Assurance Manual* (EASOPQAM, EPA 2002). Sample handling and chain-of-custody followed guidelines described in Section 2 of the EASOPQAM.

#### 2.3 SAMPLE COLLECTION

Before sampling began at each water station, *in-situ* measurements of temperature, pH, conductivity, and dissolved oxygen readings were taken at mid-depth using a YSI 6920 multi-probe sonde and recorded in a bound field book (Table 1). Global Positioning System readings were recorded at each sediment sampling location using a Garmin 76S (Garmin 2002).

A surface water sample for PAH analyses was obtained as a single grab from the middle of each fished reach before the sediment or fish samples were collected. Sediment samples consisted of a composite of three surface samples collected with a scoop and bracket from a depth of 0-5 inches. Each sediment composite consisted of: one portion taken at the beginning of each fished reach, a second from the middle of the reach and a third from the end of the reach. No sediment or fish samples were collected from Station CS-04 due to inaccessibility. GPS locations for surface water and sediment sampling are in Table 2.

Fish were collected using an electro-fishing equipped boat. The size and species of the fish obtained depended upon their availability in each reach of the slough (Table 3). An attempt was made to collect the same species at each station. After collection, fish were measured, weighed, scaled or skinned, filleted, frozen, and processed for chemical analyses in accordance with the EASOPQAM. The tissue processing was completed in the Ecological Assessment Branch (EAB) Tissue Processing Lab. After preparation the tissue samples were sent to the SESD Analytical Support Branch (ASB) for analyses according to the Analytical Support Branch Laboratory Operations and Quality Assurance Manual (ASBLOQAM 2005) (Table 3).

#### 3.0 RESULTS/DISCUSSION

#### 3.1 SURFACE WATER

The PAH concentrations were compared to the Ambient Water Quality Criteria for Human Health (AWQC-HH) by personal communication with Kevin Koporek, a Region 4, Toxicologist. The results of surface water samples from the background station (CS-01), and stations CS-02 and CS-03, were not detected at or below the minimum reporting limit (MRL) requested for PAHs. At station CS-03, there was an oil sheen present on the water surface. The surface water sample from CS-04 contained four PAHs and a strong creosote odor was noted as the sampler walked out into the slough to collect the sample. CS-04 is located in close proximity to the seep that was observed during the reconnaissance.

Detected PAHs from CS-04 are shown in Table 4. Detected PAHs were reviewed by the toxicologist whose report stated that the reported levels of acenaphthene and fluorene were well below the federal AWQC-HH. The two AWQC-HH values are based on 1) eating aquatic organisms; 2) drinking the water and ingesting aquatic organisms. There are no AWQC-HH for 2-methylnaphthalene or phenanthrene. The reported levels for these two compounds, however, were well below the AWQC-HH for other noncarcinogenic PAH compounds (EPA 2006a,b).

#### 3.2 SEDIMENTS

The PAH concentrations were compared to the Soil Preliminary Remediation Goals (PRG)/ranges. PRGs are tools for evaluating and cleaning up soils at contaminated sites (US EPA 2004). They are risk-based concentrations that are intended to assist risk assessors and others in initial screening-level evaluations of environmental measurements. The PRGs contained in the Region 9 PRG Table are generic; they are calculated without site specific information. PRGs should be viewed as Agency guidelines, not legally enforceable standards. They are used for site "screening" and as initial cleanup goals if applicable (EPA 2004).

No PAHs were detected at or below the MRL in sediment samples CS-01 and CS-02. Seven PAHs were detected in the sediment at CS-03 (See Table 4). Two tentatively identified compounds (TIC) were identified as well at CS-03. No sediment sample was collected from CS-04 because the presence of extremely soft sediments caused unsafe conditions and there was no boat access. Based on the Region 9 PRGs, the concentrations of the contaminants present in the sediment samples were below or within the EPA risk range based on human contact with soils. This is a conservative assessment for human health as the exposure to sediments should be much less than for residential soils.

#### 3.3 FISH

In all fish tissue samples, PAHs were not detected and the reporting limits were satisfactory. An EPA reference on considering fish tissue data, *Guidance for Assessing Chemical Contaminant Data for use in Fish Advisories, Vol. 1* (US EPA. 2000), states that fish have the ability to metabolize PAHs and this may explain why they are frequently not detected or found at very low concentrations even from areas that may be heavily contaminated with PAHs.

#### 4.0 CONCLUSION

Overall, levels of PAHs detected in surface water, sediment, and fish tissues during the May, 2006, sampling event are not of concern for human exposure at this time. However, EPA recommends additional sampling in the seep area, identified during the reconnaissance, to further delineate any potential problems that may be associated with the seep.

#### 5.0 REFERENCES

- GARMIN 2002. GARMIN GPSMAP 76S Chartplotting Receiver Manual and Reference Guide. GARMIN International, Inc. Olathe, Kansas.
- US EPA 2006a. National Recommended Water Quality Criteria: 2006. Office of Water/Office of Science and Technology. [http://www.epa.gov/waterscience/criteria/wqcriteria.html]
- US EPA 2006b. EPA Region III Risk Based Concentration Table for Human Health Risk Assessment, October 2006 update. [http://www.epa.gov/reg3hwmd/risk/human/index.htm]
- US EPA 2005. Analytical Support Branch Laboratory Operations and Quality Assurance Manual. US Environmental Protection Agency, Region 4, Science and Ecosystem Support Division, Analytical Support Branch, Athens, GA.
- US EPA 2004. Preliminary Remediation Goal (PRGs), EPA Region 9, updated October 2004. [http://www.epa.gov/region09/waste/sfund/prg/index.html]
- US EPA 2003. Draft Preliminary Assessment/Site Inspection Report, Gulf States Creosoting Company prepared by Weston Solutions, Inc, expressly for EPA.
- US EPA 2002. Ecological Assessment Standard Operating Procedures and Quality Assurance Manual. US Environmental Protection Agency, Region 4, Science and Ecosystem Support Division, Ecological Assessment Branch, Athens, GA.
- US EPA 2000b. Guidance for Assessing Chemical Contaminant Data for Use in Fish Advisories, Third Edition, EPA 823-B-00-007, November 2000. http://www.epa.gov/waterscience/fish/guidance.html]

Table 1: In-situ Measurements at Creosote Slough, May, 2006.

Station (Total Depth)	Date/Time 05/31/06	Sample Depth	Temperature °C	1	Dissolved Oxygen mg/L	pН
CS-01 (3.0 ft.)	0745	1.7 ft.	25.99	35	1.33	4.46
CS-02 (6.5 ft.)	0940	3.5 ft.	23.95	101	0.45	5.16
CS-03 (4.5 ft.)	1350	2.0 ft.	27.90	80	8.84	6.01
CS-04 (2.5 ft.)	1310	0.5 ft.	25.63	117	0.54	5.25

Table 2: Latitude and Longitude of Creosote Slough Sampling Stations, May, 2006

<b>G</b> F	S Coordinates*
32° 19.069' N	90° 08.963' W
32° 19.077' N	90° 08.941' W
32° 19.082' N	90° 08.961' W
22° 18 705' N	90° 08.844' W
32° 18.780' N	90° 08.890' W
32° 18.780' N	90° 08.820' W
32° 18 495' N	90° 08.952' W
32° 18.464' N	90° 09.086' W
32° 18.442' N	90° 09.231' W
220 10 621' N	90° 08.779' W
	32° 19.077' N 32° 19.082' N  32° 18.795' N 32° 18.780' N 32° 18.780' N 32° 18.495' N 32° 18.464' N

<sup>\*</sup> North American Datum (NAD) 1983.

<sup>\*\*</sup> Fish were collected along the entire A to B reach for each station.

Table 3. Physical p	Table 3. Physical parameters of fish samplescollected from Creosote Slough, May, 2006 -6-							
		-						
STATION-FISH ID	DATE	DATE		TOTAL	TOTAL	1	R. FILET WT.	
	COLLECTED	FILLETED	WT (g)	LENGTH (mm	FILET WT. (g)	(g)	(g)	ANALYSIS ID
CS01-BLC1	5/31/2006	6/2/2006	243			55	1	Composite of 2
CS01-BLC2	5/31/2006	6/2/2006	186	245	81	41	40	CS01-BLC
CS01-RSF1	5/31/2006	6/2/2006	84	160	27	. 15		Composite of 3
CS01-RSF2	5/31/2006	6/2/2006	109	165	40	21		CS01-RSF
CS01-RSF3	5/31/2006	6/2/2006	115	175	41	21	20	
CS01-LMB1	5/31/2006	6/2/2006	1248	425	403	210		Composite of 2
CS01-LMB2	5/31/2006	6/2/2006	1011	410	381	194	216	CS01-LMB
CS01-LCS1	5/31/2006	6/2/2006	385	290	134	68	66	Single
CS01-SPG1	5/31/2006	6/2/2006	877	650	223	104	119	Single
CS01-BRB1	5/31/2006	6/2/2006	990	405	282	152	130	Composite of 2
CS01-BRB2	5/31/2006	6/2/2006	890	390	264	124	141	CS01-BRB
		·				**************************************		
CS02-BLC1 (L)	5/31/2006	6/2/2006	208	235	93	48	45	Composite of 3
CS02-BLC2 (L)	5/31/2006	6/2/2006	193	230	` 84	43		Large
CS02-BLC3 (L)	5/31/2006	6/2/2006	193	225	87	46		CS02-LBCL
CS02-LMB1	5/31/2006	6/2/2006		380	321	171		Composite of 2
CS02-LMB2	5/31/2006	6/2/2006	644	360	247	129		CS02-LMB
CS02-SPG1	5/31/2006	6/2/2006	292	460	. 79	39		Single
CS02-LCS1	5/31/2006	6/2/2006	261	255	108	58		Single
CS02-BLG1	5/31/2006	6/2/2006	120	174	43	22		Composite of 5
CS02-BLG2	5/31/2006	6/2/2006	74	152	30	15		CS02-BLG
CS02-BLG3	5/31/2006	6/2/2006	92	158	37	19	18	
CS02-BLG4	5/31/2006	6/2/2006	68	152	28	14	14	
CS02-BLG5	5/31/2006	6/2/2006	68	149	26	15	11	
CS02-BLC1 (S)	5/31/2006	6/2/2006	62	162	26	14	12	Composite of 3
CS02-BLC2 (S)	5/31/2006	6/2/2006	62	162	26	14		Small
CS02-BLC3 (S)	5/31/2006	6/2/2006	51	163	22	12	10	CS02-BLCS
				• • • • • • • • • • • • • • • • • • • •				
CS03-BLC1	5/31/2006	6/2/2006	106	190	93	48		Composite of 2
CS03-BLC2	5/31/2006	6/2/2006	202	230	44	23	21	CS03-BLC
CS03-BLG1	5/31/2006	6/2/2006	64	145	26	14	12	Composite of 3
CS03-BLG2	5/31/2006	6/2/2006	58	140	21	. 11		CS03-BLG
CS03-BLG3	5/31/2006	6/2/2006	51	135	18	10	8	
CSO3-SPS1	5/31/2006	6/2/2006	964	410	383	199	184	Single
CS03-BRB1	5/31/2006	6/2/2006	759	355	231	121		Single

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Table 3. Continued							<b>-</b> 7-	_
STATION-FISH ID	DATE	DATE	TOTAL	TOTAL	TOTAL		R. FILET WT.	COMMENTS
	COLLECTED	FILLETED	WT (g)	LENGTH (mm)	FILET WT. (g)	(g)	(g)	
CS03-RDS1	5/31/2006	6/2/2006	142	190	50	26	24	Composite of 5
CS03-RDS2	5/31/2006	6/2/2006	146	189	52	27	25	CS03-RDS
CS03-RDS3	5/31/2006	6/2/2006	137	189	45	24	21	
CS03-RDS4	5/31/2006	6/2/2006	143	190		26	25	
CS03-RDS5	5/31/2006	6/2/2006	85	160	31	17	14	
CS03-LMB1	5/31/2006	6/2/2006	496	325	189	99	90	Composite of 4
CS03-LMB2	5/31/2006	6/2/2006	278	274	107	55	52	CS03-LMB
CS03-LMB3	5/31/2006	6/2/2006	432	314	178	92	. 86	
CS03-LMB4	5/31/2006	6/2/2006	430	314	178	90	88	
								-
BLC = BLACK CRAPE			IS					
BLG = BLUEGILL - <i>Le</i>								
BRB = BROWN BULL			S					•
LCS = LAKE CHUB SI	JCKER - Erimy	zon sucetta						
LMB = LARGEMOUTH								
RDS = REDBREAST S								
RSF = REDEAR SUNFISH - Lepomis microlophus								
SPG = SPOTTED GAR - Lepisosteus oculatus								
SPS = SPOTTED SUC	CKER - Minytrer	na melanops	;					

4 5 1

Table 4: Detected PAHs in surface water and sediment in Creosote Slough, May, 2006.

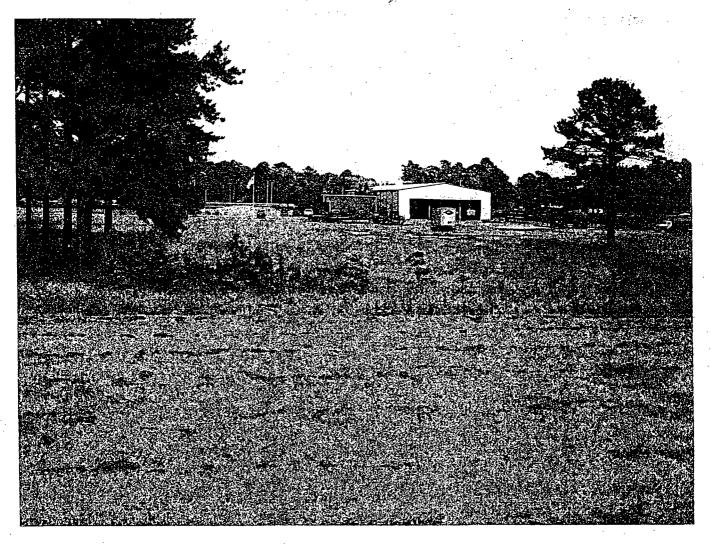
Surface Water (CS-04) Grab	μg/L	AWQC* μg/L
2-Methylnaphthalene	1.8	NA
Acenaphthene	12.0	670/990
Fluorene	9.3	1100/5300
Phenanthrene	2.8	NA
Sediments (CS-03) Composite	mg/kg	Soil PRG/range mg/kg
Fluoranthene	0.2	2300
Pyrene	0.13	2300
Benzo(a)anthracene	0.078	0.62-62
Chrysene	0.13	62-6200
Benzo(b)fluoranthene	0.12	0.62-62
Benzo(k)fluoranthene	0.12	6.2-620
Benzo(a)pyrene	0.066	0.062-6.2

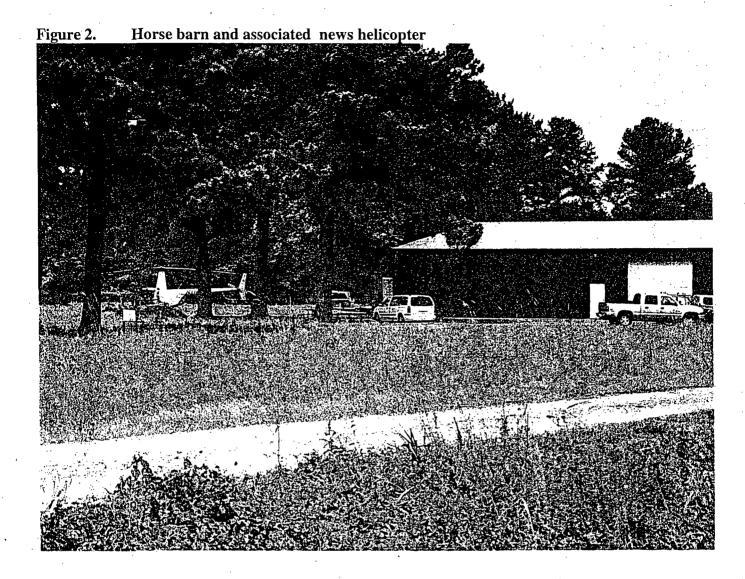
AWQC = Ambient Water Quality Criteria

\* Chronic/Acute water quality criteria
PRG = Preliminary Remediation Goal



Figure 1. ConSteel Co., Inc





Sept.

Service B

Figure 3. Levee separating former Gulf States Creosoting from Creosote Slough

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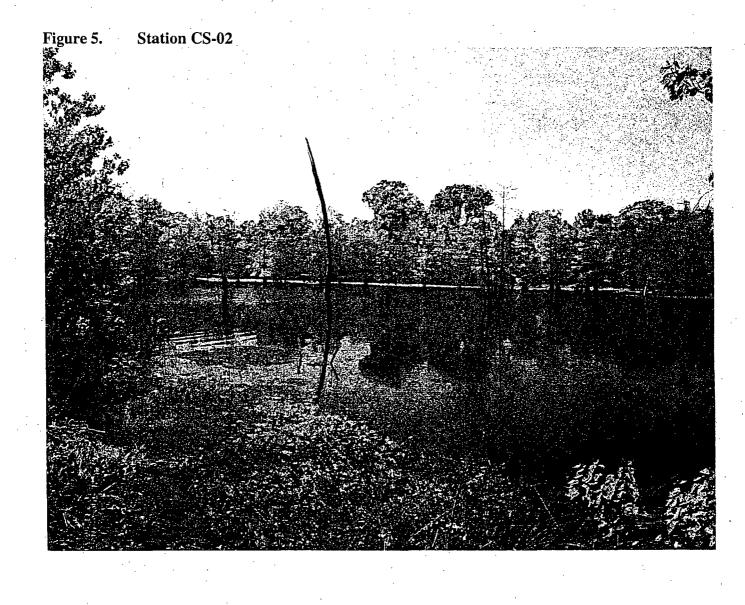
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Figure 4. CS-01 Reference Station





(A)

Figure 6. Stations CS-03 and CS-04

How we



Figure 7. Seep Area

12 11 25

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# APPENDIX I

Water, Sediment and Tissue Data

#### **EXTRACTABLES SAMPLE ANALYSIS**

EPA - REGION IV SESD, ATHENS, GA

Production Date: 08/16/2006 14:39

Sample 5727 FY 2006 Project: 06-0436

PAH Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS-01-SW / Media: SURFACE WATER Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006 07:45

Ending:

RESULTS	UNITS	ANALYTE
1.2 U	UG/L	2-Methylnaphthalene
1.4 U	UG/L	Naphthalene
1.3 U	UG/L	Acenaphthylene
1.2 U	UG/L	Acenaphthene
1.5 U	UG/L	Fluorene
1.5 U	UG/L	Phenanthrene
1.6 U	UG/L	Anthracene
1.6 U	UG/L	Fluoranthene
1.5 U	UG/L	Pyrene
1.3 U	UG/L	Benzo(a)Anthracene
1.5 U	UG/L	Chrysene
1.0 U	UG/L	Benzo(b)Fluoranthene
1.5 U	UG/L	. Benzo(k)Fluoranthene
1:0 U	UG/L	Benzo-a-Pyrene
1.2 U	UG/L	Indeno (1,2,3-cd) Pyrene
1.3 U	UG/L	Dibenzo(a,h)Anthracene
1.0 U	UG/L	Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification.

K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### EXTRACTABLES SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

Sample 5728 FY 2006

06 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS-01-SD /
Media: SEDIMENT

Negative

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006 08:00

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
39 UJ	UG/KG	2-Methylnaphthalene
39 U	UG/KG	Naphthalene
39 U	UG/KG	Acenaphthylene
39 U	UG/KG	Acenaphthene
39 U	UG/KG	Fluorene
39 U	UG/KG	Phenanthrene
39 U	UG/KG	Anthracene
39 U	UG/KG	Fluoranthene
39 U	UG/KG	Pyrene
39 U	UG/KG	Benzo(a)Anthracene
39 U	UG/KG	Chrysene
39 U	UG/KG	Benzo(b)Fluoranthene
39 U	UG/KG	Benzo(k)Fluoranthene
39 U	UG/KG	Benzo-a-Pyrene
39 U	UG/KG	Indeno (1,2,3-cd) Pyrene
39 U	UG/KG	Dibenzo(a,h)Anthracene
58	%	% Moisture
39 U	UG/KG	Benzo(ghi)Perylene

J-qualified: 2-Methylnaphthalene recovery outside QC Limits

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### **EXTRACTABLES SAMPLE ANALYSIS EPA - REGION IV SESD, ATHENS, GA** Production Date: 08/16/2006 14:39 Produced by: Revell, Dennis Sample 5729 FY 2006 Project: 06-0436 Requestor: Donna Webster **PAH Scan** Project Leader: PMEYER Facility: Gulf States Creosoting Flowood, MS Beginning: 05/31/2006 09:40 Program: SF Ending: Id/Station: CS-02-SW / Media: SURFACE WATER **RESULTS UNITS ANALYTE** 1.1 U UG/L 2-Methylnaphthalene 1.3 U UG/L Naphthalene 1.2 U UG/L Acenaphthylene 1.1 U UG/L Acenaphthene 1.4 U UG/L Fluorene 1.4 U UG/L Phenanthrene 1.5 U UG/L Anthracene 1.5 U UG/L Fluoranthene 1.4 U UG/L Pyrene 1.2 U UG/L Benzo(a)Anthracene

1.4 U

1.0 U

1.4 U

1.0 U

1.1 U

1.2 U

1.0 U

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

Chrysene

Benzo(b)Fluoranthene

Benzo(k)Fluoranthene

Indeno (1,2,3-cd) Pyrene

Dibenzo(a,h)Anthracene

Benzo(ghi)Perylene

Benzo-a-Pyrene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### **EXTRACTABLES SAMPLE ANALYSIS EPA - REGION IV SESD, ATHENS, GA** Production Date: 08/16/2006 14:39 Produced by: Revell, Dennis Sample 5730 FY 2006 Project: 06-0436 Requestor: Donna Webster **PAH Scan** Project Leader: PMEYER Facility: Gulf States Creosoting Flowood, MS Beginning: 05/31/2006 09:55 Program: SF Endina: Id/Station: CS-02-SD / Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS RESULTS UNITS **ANALYTE** 72 UJ UG/KG 2-Methylnaphthalene 72 U UG/KG Naphthalene 72 U UG/KG Acenaphthylene 72 U UG/KG Acenaphthene 72 U UG/KG Fluorene 72 U UG/KG Phenanthrene 72 U UG/KG Anthracene 72 U Fluoranthene UG/KG 72 U UG/KG Pyrene 72 U UG/KG Benzo(a)Anthracene 72 U UG/KG Chrysene 72 U UG/KG Benzo(b)Fluoranthene

J-qualified: 2-Methylnaphthalene recovery outside QC Limits

72 U

72 U

72 U

72 U

72 U

77

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

Benzo(k)Fluoranthene

Indeno (1,2,3-cd) Pyrene

Dibenzo(a.h)Anthracene

Benzo(ghi)Pervlene

Benzo-a-Pyrene

% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be less than the reported value L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### XTRACTABLES SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

Sample 5732 FY 2006 Project: 06-0436

PAH Scan

Project: 05-043t

Flowood, MS

Produced by: Revell, Dennis Requestor: Donna Webster

Requestor: Donna Webster Project Leader: PMEYER

Beginning: 05/31/2006 13:50

Ending:

Facility: Gulf States Creosoting Program: SF

0.95 U

UG/L

Id/Station: CS-03-SW / Media: SURFACE WATER

RESULTS UNITS **ANALYTE** 1.2 U UG/L 2-Methylnaphthalene 1.2 U UG/L Naphthalene 1.1 U UG/L Acenaphthylene 1.0 U UG/L Acenaphthene 1.3 U UG/L Fluorene 1.3 U UG/L Phenanthrene 1.4 U UG/L Anthracene 1.4 U UG/L Fluoranthene 1.3 U UG/L Pyrene 1.1 U UG/L Benzo(a)Anthracene 1.3 U UG/L Chrysene 0.95 U UG/L Benzo(b)Fluoranthene 1.3 U Benzo(k)Fluoranthene UG/L 0.95 U Benzo-a-Pyrene UG/L 1.0 U UG/L Indeno (1,2,3-cd) Pyrene 1.1 U UG/L Dibenzo(a,h)Anthracene

Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

#### XTRACTABLES SAMPLE ANALYSIS

#### **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/16/2006 14:39

Sample **PAH Scan** 

5733 FY 2006 Project: 06-0436

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

id/Station: CS-03-SD / Media: SEDIMENT

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006 14:00

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
61 UJ	UG/KG	2-Methylnaphthalene
61 U	UG/KG	Naphthalene
61°U	UG/KG	Acenaphthylene
61 U	UG/KG	Acenaphthene
61 U	UG/KG	Fluorene
61 U	UG/KG	Phenanthrene
61 U	UG/KG	Anthracene
200	UG/KG	Fluoranthene
130	UG/KG	Pyrene
78	UG/KG	Benzo(a)Anthracene
130	UG/KG	Chrysene
120	UG/KG	Benzo(b)Fluoranthene
120	UG/KG	Benzo(k)Fluoranthene
66	UG/KG	Benzo-a-Pyrene
61 U	UG/KG	Indeno (1,2,3-cd) Pyrene
61 U	UG/KG	Dibenzo(a,h)Anthracene
73	<b>%</b> ,	% Moisture
61 U	UG/KG	Benzo(ghi)Perylene

J-qualified: 2-Methylnaphthalene recovery outside QC Limits

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### XTRACTABLES SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

Sample 5731 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006 13:10

Produced by: Revell, Dennis

Ending:

Program: SF

Id/Station: CS-04-SW / Media: SURFACE WATER

RESULTS	UNITS	ANALYTE
1.8	UG/L	2-Methylnaphthalene
1.3 U	UG/L	Naphthalene
1.2 U	UG/L	Acenaphthylene
12	UG/L	Acenaphthene
9.3	UG/L	Fluorene
2.8	UG/L	Phenanthrene
1.5 U	UG/L	Anthracene
1.5 U	UG/L	Fluoranthene
1.4 U	UG/L	Pyrene
1.2 U	UG/L	Benzo(a)Anthracene
1.4 U	UG/L	Chrysene
. 1.0 U	UG/L	Benzo(b)Fluoranthene
1.4 U	UG/L	Benzo(k)Fluoranthene
1.0 U	UG/L	Benzo-a-Pyrene
1.1 U	UG/L	Indeno (1,2,3-cd) Pyrene
1.2 U	UG/L	Dibenzo(a,h)Anthracene
1.0 U	UG/L	Benzo(ghi)Perylene
		, .,

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### **EXTRACTABLES SAMPLE ANALYSIS**

EPA - REGION IV SESD, ATHENS, GA

Production Date: 08/28/2006 10:39

Sample 6082 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS01-BLC /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS UNITS **ANALYTE** 0.016 U MG/KG 2-Methylnaphthalene 0.016 U MG/KG Naphthalene 0.016 U MG/KG Acenaphthylene 0.016 U MG/KG Acenaphthene 0.016 U MG/KG Fluorene 0.016 U MG/KG Phenanthrene 0.016 U MG/KG Anthracene 0.016 U MG/KG Fluoranthene 0.016 U MG/KG Pyrene 0.016 U MG/KG Benzo(a)Anthracene 0.016 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.016 U MG/KG Benzo(k)Fluoranthene 0.016 U MG/KG Benzo-a-Pyrene 0.018 U MG/KG Indeno (1,2,3-cd) Pyrene 0.018 U MG/KG Dibenzo(a,h)Anthracene 0.016 U MG/KG Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte reported as tentative identification. | Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be less than the reported value L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/28/2006 10:39

Sample 6086 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Media: FISH

Id/Station: CS01-RSF /

Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Produced by: Revell, Dennis

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.017 U	MG/KG	2-Methylnaphthalene
0.017 U	MG/KG	Naphthalene
0.017 U	MG/KG	Acenaphthylene
0.017 U	MG/KG	Acenaphthene
0.017 U	MG/KG	Fluorene
0.017 U	MG/KG	Phenanthrene
0.017 U	MG/KG	Anthracene
0.017 U	MG/KG	Fluoranthene
0.017 U	MG/KG	Pyrene
0.017 U	MG/KG	Benzo(a)Anthracene
0.017 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.017 U	MG/KG	Benzo(k)Fluoranthene
0.017 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0,018 U	MG/KG	Dibenzo(a,h)Anthracene
0.017 U	MG/KG	Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Sample

6085 FY 2006

Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS01-LMB /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.017 U	MG/KG	2-Methylnaphthalene
0.017 U	MG/KG	Naphthalene
0.017 U	MG/KG	Acenaphthylene
0.017 U	MG/KG	Acenaphthene
0.017 U	MG/KG	Fluorene
0.017 U	MG/KG	Phenanthrene
0.017 U	MG/KG	Anthracene
0.017 U	MG/KG	Fluoranthene
0.017 U	MG/KG	Pyrene
0.017 U	MG/KG	Benzo(a)Anthracene
0.017 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.017 Ù	MG/KG	Benzo(k)Fluoranthene
0.017 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.017 U	MG/KG	Benzo(ghi)Perylene
•		•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

EPA - REGION IV SESD, ATHENS, GA

Production Date: 08/28/2006 10:39

Sample 6084 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS01-LCS1 /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Endina:

DATA REPORTED ON WET WEIGHT BASIS

**RESULTS UNITS ANALYTE** 0.017 U MG/KG 2-Methylnaphthalene 0.017 U MG/KG Naphthalene 0.017 U MG/KG Acenaphthylene 0.017 U MG/KG Acenaphthene 0.017 U MG/KG Fluorene 0.017 U MG/KG Phenanthrene 0.017 U MG/KG Anthracene 0.017 U MG/KG Fluoranthene 0.017 U MG/KG Pyrene 0.017 U MG/KG Benzo(a)Anthracene 0.017 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.017 U MG/KG Benzo(k)Fluoranthene 0.017 U MG/KG Benzo-a-Pyrene 0.018 U MG/KG Indeno (1,2,3-cd) Pyrene 0.018 U MG/KG Dibenzo(a.h)Anthracene 0.017 U MG/KG Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/28/2006 10:39

Sample 6087 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: CS01-SPG1 /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.016 U	MG/KG	2-Methylnaphthalene
0.016 U	MG/KG	Naphthalene
0.016 U	MG/KG	Acenaphthylene
0.016 U	MG/KG	Acenaphthene
0.016 U	MG/KG	Fluorene
0.016 U	MG/KG	Phenanthrene
0.016 U	MG/KG	Anthracene
0.016 U	MG/KG	Fluoranthene
0.016 U	MG/KG	Pyrene
0.016 U	MG/KG	Benzo(a)Anthracene
0.016 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.016 U	MG/KG	Benzo(k)Fluoranthene
0.016 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.016 U	MG/KG	Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

RACTABLES SAMPLE ANALYSIS EPA - REGION IV SESD, ATHENS, GA

Production Date: 11/15/2006 14:10

Sample 6090 FY 2006 Project: 06-0436

PAH Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS02-BLG /

DECILITE LINUTO

Media: FISH

Produced by: Revell, Dennis

Requestor: Donna Webster Project Leader: PMEYER

Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

KESUL15	UNITS	ANALYTE
0.016 U	MG/KG	2-Methylnaphthalene
0.016 U	MG/KG	Naphthalene
0.016 U	MG/KG	Acenaphthylene
0.016 U	MG/KG	Acenaphihene
0.016 U	MG/KG	Fluorene
0.016 U	MG/KG	Phenanthrene '
0.016 U	MG/KG	Anthracene
0.016 U	MG/KG	Fluoranthene
0.016 U	MG/KG	Pyrene
0.016 U	MG/KG	Benzo(a)Anthracene
0.016 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.016 U	MG/KG	Benzo(k)Fluoranthene
0.016 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.016 U	MG/KG	Benzo(ghi)Perylene

Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

dentification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

Sample 5729 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS-02-SW / Media: SURFACE WATER

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006 09:40

Ending:

UNITS	ANALYTE
UG/L	2-Methylnaphthalene
UG/L	Naphthalene
UG/L	Acenaphthylene
UG/L	Acenaphthene
UG/L	Fluorene
UG/L	Phenanthrene
UG/L	Anthracene
UG/L	Fluoranthene
UG/L	Pyrene .
UG/L	Benzo(a)Anthracene
UG/L	Chrysene
UG/L	Benzo(b)Fluoranthene
UG/L	Benzo(k)Fluoranthene
UG/L	Benzo-a-Pyrene
UG/L	Indeno (1,2,3-cd) Pyrene
UG/L	Dibenzo(a,h)Anthracene
UG/L	Benzo(ghi)Perylene
	UG/L UG/L UG/L UG/L UG/L UG/L UG/L UG/L

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

Sample 5730 FY 2006

Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

72 U

72 U

77

UG/KG

UG/KG

ld/Station: CS-02-SD / Media: SEDIMENT

Produced by: Revell, Dennis

Requestor: Donna Webster

Beginning: 05/31/2006 09:55

Project Leader: PMEYER

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

**ANALYTE** 

**RESULTS UNITS** 72 UJ UG/KG 2-Methylnaphthalene 72 U UG/KG Naphthalene 72 U UG/KG Acenaphthylene 72 U UG/KG Acenaphthene 72 U UG/KG Fluorene 72 U UG/KG Phenanthrene 72 U UG/KG Anthracene 72 U UG/KG Fluoranthene 72 U UG/KG Pyrene 72 U UG/KG Benzo(a)Anthracene 72 U UG/KG Chrysene 72 U UG/KG Benzo(b)Fluoranthene 72 U Benzo(k)Fluoranthene UG/KG 72 U UG/KG Benzo-a-Pyrene 72 U UG/KG Indeno (1,2,3-cd) Pyrene

Dibenzo(a.h)Anthracene

Benzo(ghi)Pervlene

% Moisture

J-qualified: 2-Methylnaphthalene recovery outside QC Limits

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39 -

Sample **PAH Scan** 

5732 FY 2006

Project: 06-0436

Facility: Gulf States Creosoting

Program: SF

1.3 U

1.0 U

1.1 U

0.95 U

0.95 U

UG/L

UG/L

UG/L

UG/L

UG/L

Id/Station: CS-03-SW / Media: SURFACE WATER Flowood, MS

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006 13:50 Ending:

**RESULTS UNITS ANALYTE** 1.2 U UG/L 2-Methylnaphthalene 1.2 U UG/L Naphthalene 1.1 U UG/L Acenaphthylene 1.0 U UG/L Acenaphthene 1.3 U UG/L Fluorene 1.3 U UG/L Phenanthrene 1.4 U UG/L Anthracene 1.4 U UG/L Fluoranthene 1.3 U UG/L Pyrene 1.1 U UG/L Benzo(a)Anthracene 1.3 U UG/L Chrysene 0.95 U UG/L Benzo(b)Fluoranthene

Benzo(k)Fluoranthene

Indeno (1,2,3-cd) Pyrene

Dibenzo(a,h)Anthracene

Benzo(ghi)Perylene

Benzo-a-Pyrene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Facility: Gulf States Creosoting

#### **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/16/2006 14:39

Sample 5733 FY 2006

Project: 06-0436

Flowood, MS

Produced by: Revell, Dennis

Requestor: Donna Webster Project Leader: PMEYER

Beginning: 05/31/2006 14:00

Ending:

Program: SF Id/Station: CS-03-SD /

Media: SEDIMENT

**PAH Scan** 

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS UNITS **ANALYTE** 61 UJ 2-Methylnaphthalene UG/KG 61 U UG/KG Naphthalene 61 U UG/KG Acenaphthylene 61 U UG/KG Acenaphthene 61 U UG/KG Fluorene 61 U UG/KG Phenanthrene 61 U UG/KG Anthracene 200 UG/KG Fluoranthene 130 UG/KG Pyrene 78 UG/KG Benzo(a)Anthracene 130 UG/KG Chrysene 120 UG/KG Benzo(b)Fluoranthene 120 UG/KG Benzo(k)Fluoranthene 66 UG/KG Benzo-a-Pyrene 61 U UG/KG Indeno (1,2,3-cd) Pyrene 61 U UG/KG Dibenzo(a.h)Anthracene 73 % Moisture 61 U UG/KG Benzo(ghi)Perylene

I-qualified: 2-Methylnaphthalene recovery outside QC Limits

J-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. V-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. <-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.</p>

<sup>--</sup>Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

Sample 5731 FY 2006

Project: 06-0436

**PAH Scan** Facility: Gulf States Creosoting

Program: SF

Id/Station: CS-04-SW / Media: SURFACE WATER Flowood, MS

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER

Beginning: 05/31/2006 13:10

Ending:

ESULTS	UNITS	ANALYTE
1.8	UG/L	2-Methylnaphthalene
1.3 U	UG/L	Naphthalene
1.2 U	UG/L	Acenaphthylene
12	.UG/L	Acenaphthene
9.3	UG/L	Fluorene
2.8	UG/L	Phenanthrene
1.5 U	UG/L	Anthracene
1.5 U	UG/L	Fluoranthene
1.4 U	UG/L	Pyrene
1.2 U	UG/L	Benzo(a)Anthracene
1.4 U	UG/L	Chrysene
1.0 U	UG/L	Benzo(b)Fluoranthene
1.4 U	UG/L	Benzo(k)Fluoranthene
1.0 U	UG/L	Benzo-a-Pyrene
1.1 U	UG/L	Indeno (1,2,3-cd) Pyrene
1.2 U	UG/L	Dibenzo(a,h)Anthracene
1.0 U	UG/L	Benzo(ghi)Pervlene

J-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. 4-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Sample 6082 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS01-BLC /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

**RESULTS UNITS** ANALYTE 0.016 U MG/KG 2-Methylnaphthalene 0.016 U MG/KG Naphthalene 0.016 U MG/KG Acenaphthylene 0.016 U MG/KG Acenaphthene 0.016 U MG/KG Fluorene 0.016 U MG/KG Phenanthrene 0.016 U MG/KG Anthracene 0.016 U MG/KG Fluoranthene 0.016 U MG/KG Pyrene 0.016 U MG/KG Benzo(a)Anthracene 0.016 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.016 U MG/KG Benzo(k)Fluoranthene 0.016 U MG/KG Benzo-a-Pyrene 0.018 U MG/KG Indeno (1.2.3-cd) Pyrene 0.018 U MG/KG Dibenzo(a,h)Anthracene 0.016 U MG/KG Benzo(ghi)Perylene

J-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is p

<sup>-</sup>Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

VA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

<sup>3-</sup>Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Sample

6086 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS01-RSF /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.017 U	MG/KG	2-Methylnaphthalene
0.017 U	MG/KG	Naphthalene
0.017 U	MG/KG	Acenaphthylene
0.017 U	MG/KG	Acenaphthene
0.017 U	MG/KG	Fluorene
0.017 U	MG/KG	Phenanthrene
0.017 U	MG/KG	Anthracene .
0.017 U	MG/KG	Fluoranthene
0.017 U	MG/KG	Pyrene
0.017-U	MG/KG	Benzo(a)Anthracene
0.017 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.017 U	MG/KG	Benzo(k)Fluoranthene
0.017 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.017 U	MG/KG	Benzo(ghi)Perylene

J-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD. ATHENS. GA** 

Production Date: 08/28/2006 10:39

Sample FY 2006 6085 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

MG/KG

MG/KG

MG/KG

Flowood, MS

Program: SF

Id/Station: CS01-LMB /

Media: FISH

0.018 U

0.018 U

0.017 U

DATA REPORTED ON WET WEIGHT BASIS

Produced by: Revell, Dennis

Requestor: Donna Webster

Project Leader: PMEYER

Beginning: 05/31/2006

Endina:

**RESULTS UNITS** ANALYTE 0.017 U MG/KG 2-Methylnaphthalene 0.017 U MG/KG Naphthalene 0.017 U MG/KG Acenaphthylene 0.017 U MG/KG Acenaphthene 0.017 U MG/KG Fluorene 0.017 U MG/KG Phenanthrene 0.017 U MG/KG Anthracene 0.017 U MG/KG Fluoranthene 0.017 U MG/KG Pyrene 0.017 U MG/KG Benzo(a)Anthracene 0.017 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.017 U MG/KG Benzo(k)Fluoranthene 0.017 U MG/KG Benzo-a-Pyrene

Indeno (1,2,3-cd) Pyrene

Dibenzo(a,h)Anthracene

Benzo(ghi)Perylene

J-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

<sup>--</sup>Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate, Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Sample 6084 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Media: FISH

Id/Station: CS01-LCS1 /

Requestor: Donna Webster Project Leader: PMEYER

Beginning: 05/31/2006

Produced by: Revell, Dennis

Ending:

DATA REPORTED ON WET WEIGHT BASIS

**RESULTS UNITS** ANALYTE 0.017 U MG/KG 2-Methylnaphthalene 0.017 U MG/KG Naphthalene 0.017 U MG/KG Acenaphthylene Acenaphthene 0.017 U MG/KG 0.017 U MG/KG Fluorene 0.017 U MG/KG Phenanthrene 0.017 U MG/KG Anthracene 0.017 U MG/KG Fluoranthene 0.017 U MG/KG Pyrene 0.017 U MG/KG Benzo(a)Anthracene 0.017 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.017 U MG/KG Benzo(k)Fluoranthene 0.017 U MG/KG Benzo-a-Pyrene 0.018 U MG/KG Indeno (1,2,3-cd) Pyrene 0.018 U MG/KG Dibenzo(a,h)Anthracene 0.017 U MG/KG Benzo(ghi)Pervlene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

EPA - REGION IV SESD, ATHENS, GA

Production Date: 08/28/2006 10:39

Sample 6083 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS01-BRB /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.016 U	MG/KG	2-Methylnaphthalene
0.016 U	MG/KG	Naphthalene
0.016 U	MG/KG	Acenaphthylene
0.016 U	MG/KG	Acenaphthene
0.016 U	MG/KG	Fluorene
0.016 U	MG/KG	Phenanthrene
0.016 U	MG/KG	Anthracene
0.016 U	MG/KG	Fluoranthene
0.016 U	MG/KG	Pyrene
0.016 U	MG/KG	Benzo(a)Anthracene
0.016 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.016 U	MG/KG	Benzo(k)Fluoranthene
0.016 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.016 U	MG/KG	Benzo(ghi)Perylene
		•

J-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. | V-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | C-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

<sup>--</sup>Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

VA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Sample 6088 FY 2006

Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS02-BLCL /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster

Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.017 U	MG/KG	2-Methylnaphthalene
0.017 U	MG/KG	Naphthalene
0.017 U	MG/KG	Acenaphthylene
0.017 U	MG/KG	Acenaphthene
0.017 U	MG/KG	Fluorene
0.017 U	MG/KG	Phenanthrene
0.017 U	MG/KG	Anthracene
0.017 U	MG/KG	Fluoranthene
0.017 U	MG/KG	Pyrene
0.017 U	MG/KG	Benzo(a)Anthracene
0.017 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.017 U	MG/KG	Benzo(k)Fluoranthene
0.017 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.017 U	MG/KG	Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/28/2006 10:39

Sample Project: 06-0436 6092 FY 2006

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

**PAH Scan** 

Id/Station: CS02-LMB /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

	RESULTS	UNITS	ANALYTE
	0.016 U	MG/KG	2-Methylnaphthalene
	0.016 U	MG/KG	Naphthalene
	0.016 U	MG/KG	Acenaphthylene
	0.016 U	MG/KG	Acenaphthene
	0.016 U	MG/KG	Fluorene
	0.016 U	MG/KG	Phenanthrene
	0.016 U	MG/KG	Anthracene
	0.016 U	MG/KG	Fluoranthene
	0.016 U	MG/KG	Pyrene
	0.016 U	MG/KG	Benzo(a)Anthracene
	0.016 U	MG/KG	Chrysene
-	0.019 U	MG/KG	Benzo(b)Fluoranthene
	0.016 U	MG/KG	Benzo(k)Fluoranthene
•	0.016 U	MG/KG	Benzo-a-Pyrene
	0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
	0.018 U	MG/KG	Dibenzo(a,h)Anthracene
	0.016 U	MG/KG	Benzo(ghi)Perylene
			•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Page 1 of 1

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Project: 06-0436 Sample 6093 FY 2006

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS02SPG1 /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.017 U	MG/KG	2-Methylnaphthalene
0.017 U	MG/KG	Naphthalene
0.017 U	MG/KG	Acenaphthylene
0.017 U	MG/KG	Acenaphthene
0.017 U	MG/KG	Fluorene
0.017 U	MG/KG	Phenanthrene
0.017 U	MG/KG	Anthracene
0.017 U	MG/KG	Fluoranthene
0.017 U	MG/KG	Pyrene ·
0.017 U	MG/KG	Benzo(a)Anthracene
0.017 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.017 U	MG/KG	Benzo(k)Fluoranthene
0.017 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyreno
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.017 U	MG/KG	Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Ideritification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

REGION IV SESD, ATHENS, GA

Production Date: 11/15/2006 14:10

Project: 06-0436 Sample 6091 FY 2006

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS02-LCS1 /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.016 U	MG/KG	2-Methylnaphthalene
0.016 U	MG/KG	Naphthalene
0.016 U	MG/KG	Acenaphthylene
0.016 U	MG/KG	Acenaphthene
0.016 U	MG/KG	Fluorene
0.016 U	MG/KG	Phenanthrene
0.016 U	MG/KG	Anthracene
0.016 U	MG/KG	Fluoranthene
0.016 U	MG/KG	Pyrene
0.016 U	MG/KG	Benzo(a)Anthracene
0.016 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.016 U	MG/KG	Benzo(k)Fluoranthene
0.016 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.016 U	MG/KG	Benzo(ghi)Perylene
_		

alyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. esumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. Intification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. intification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. esence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. RECONITOR CONTROL CONT

Production Date: 1775/2000 14:10

Sample 6089 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS02-BLCS /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster

Project Leader: PMEYER
Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.017 U	MG/KG	2-Methylnaphthalene
0.017 U	MG/KG	Naphthalene
0.017 U	MG/KG	Acenaphthylene
0.017 U	MG/KG	Acenaphthene
0.017 U	MG/KG	Fluorene
0.017 U	MG/KG	Phenanthrene
0.017 U	MG/KG	Anthracene
0.017 U	MG/KG	Fluoranthene
0.017 U	MG/KG	Pyrene
0.017 U	MG/KG	Benzo(a)Anthracene
0.017 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.017 U	MG/KG	Benzo(k)Fluoranthene
0.017 U	MG/KG	Benzo-a-Pyrene
0.018 ป	MG/KG	indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.017 U	MG/KG	Benzo(ghi)Perylene

nalyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. resumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. entification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

entification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Facility: Gulf States Creosoting

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Sample **PAH Scan** 

Program: SF

6094 FY 2006 Project: 06-0436

Flowood, MS

Produced by: Revell, Dennis

Requestor: Donna Webster

Project Leader: PMEYER

Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

Id/Station: CS03-BLC / Media: FISH **RESULTS UNITS ANALYTE** 0.016 U MG/KG 2-Methylnaphthalene 0.016 U MG/KG Naphthalene 0.016 U MG/KG Acenaphthylene 0.016 U MG/KG Acenaphthene

0.016 U MG/KG Fluorene 0.016 U MG/KG Phenanthrene 0.016 U MG/KG Anthracene 0.016 U MG/KG Fluoranthene

0.016 U MG/KG Pyrene 0.016 U MG/KG Benzo(a)Anthracene

0.016 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene

0.016 U MG/KG Benzo(k)Fluoranthene 0.016 U MG/KG Benzo-a-Pyrene 0.018 U MG/KG Indeno (1,2,3-cd) Pyrene

0.018 U MG/KG Dibenzo(a,h)Anthracene 0.016 U MG/KG Benzo(ghi)Perylene

J-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. V-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. C-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Facility: Gulf States Creosoting

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/28/2006 10:39

Sample 6095 FY 2006 Project: 06-0436

PAH Scan

Flowood, MS

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER

Beginning: 05/31/2006

Ending:

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Id/Station: CS03-BLG /

Media: FISH

Program: SF

DATA REPORTED ON WET WEIGHT BASIS

**RESULTS UNITS ANALYTE** 0.016 U MG/KG 2-Methylnaphthalene 0.016 U MG/KG Naphthalene 0.016 U MG/KG Acenaphthylene 0.016 U MG/KG Acenaphthene 0.016 U MG/KG Fluorene 0.016 U MG/KG Phenanthrene 0.016 U MG/KG Anthracene 0.016 U MG/KG Fluoranthene 0.016 U MG/KG Pyrene 0.016 U MG/KG Benzo(a)Anthracene 0.016 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.016 U MG/KG Benzo(k)Fluoranthene 0.016 U MG/KG Benzo-a-Pyrene Indeno (1,2,3-cd) Pyrene 0.018 U MG/KG 0.018 U MG/KG Dibenzo(a,h)Anthracene 0.016 U MG/KG Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

### EPA - REGION IV SESD, ATHENS, GA

Production Date: 08/28/2006 10:39

Sample 6099 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS03-SPS /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

**RESULTS UNITS ANALYTE** 0.016 U MG/KG 2-Methylnaphthalene 0.016 U MG/KG Naphthalene 0.016 U MG/KG Acenaphthylene 0.016 U MG/KG Acenaphthene 0.016 U MG/KG Fluorene 0.016 U MG/KG Phenanthrene 0.016 U MG/KG Anthracene 0.016 U MG/KG Fluoranthene 0.016 U MG/KG Pyrene 0.016 U MG/KG Benzo(a)Anthracene 0.016 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.016 U MG/KG Benzo(k)Fluoranthene 0.016 U MG/KG Benzo-a-Pyrene 0.018 U MG/KG Indeno (1,2,3-cd) Pyrene 0.018 U MG/KG Dibenzo(a,h)Anthracene 0.016 U MG/KG Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

FY 2006

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/28/2006 10:39

Sample 6096
PAH Scan

Facility: Gulf States Creosoting

MG/KG

MG/KG

MG/KG

.

Project: 06-0436

Program: SF

0.018 U

0.018 U

0.016 U

Id/Station: CS03-BRB /

Media: FISH

Flowood, MS

Project Leader: PMEYER Beginning: 05/31/2006 Ending:

Produced by: Revell, Dennis

Requestor: Donna Webster

DATA REPORTED ON WET WEIGHT BASIS

**RESULTS UNITS** ANALYTE 0.016 U MG/KG 2-Methylnaphthalene 0.016 U MG/KG Naphthalene 0.016 U MG/KG Acenaphthylene 0.016 U MG/KG Acenaphthene 0.016 U MG/KG Fluorene 0.016 U MG/KG Phenanthrene 0.016 U MG/KG Anthracene 0.016 U MG/KG Fluoranthene 0.016 U MG/KG Pyrene 0.016 U MG/KG Benzo(a)Anthracene 0.016 U MG/KG Chrysene 0.020 U MG/KG Benzo(b)Fluoranthene 0.016 U MG/KG Benzo(k)Fluoranthene 0.016 U MG/KG Benzo-a-Pyrene

Indeno (1,2,3-cd) Pyrene

Dibenzo(a.h)Anthracene

Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/28/2006 10:39

Sample 6098 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS03-RDS /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER

Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

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	RESULTS	UNITS	ANALYTE
	0.016 U	MG/KG	2-Methylnaphthalene
	0.016 U	MG/KG	Naphthalene
	0.016 U	MG/KG	Acenaphthylene
	0.016 U	MG/KG	Acenaphthene
	0.016 U	MG/KG	Fluorene
	0.016 U	MG/KG	Phenanthrene
	0.016 U	MG/KG	Anthracene
	0.016 U	MG/KG	Fluoranthene
r	0.016 U	MG/KG	Pyrene
	0.016 U	MG/KG	Benzo(a)Anthracene
	0.016 U	MG/KG	Chrysene
	0.020 U	MG/KG	Benzo(b)Fluoranthene
	0.016 U	MG/KG	Benzo(k)Fluoranthene
	0.016 U	MG/KG	Benzo-a-Pyrene
	0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
	0.018 U	MG/KG	Dibenzo(a,h)Anthracene
	0.016 U	MG/KG	Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

EPA - REGION IV SESD, ATHENS, GA

Production Date: 08/28/2006 10:39

Sample 6097 FY 2006 Project: 06-0436

**PAH Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: CS03-LMB /

Media: FISH

Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 05/31/2006

Ending:

DATA REPORTED ON WET WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.016 U	MG/KG	2-Methylnaphthalene
0.016 U	MG/KG	Naphthalene
0.016 U	MG/KG	Acenaphthylene
0.016 U	MG/KĠ	Acenaphthene
0.016 U	MG/KG	Fluorene
0.016 U	MG/KG	Phenanthrene
0.016 U	MG/KG	Anthracene
0.016 U	MG/KG	Fluoranthene
0.016 U	MG/KG	Pyrene
0.016 U	MG/KG	Benzo(a)Anthracene
0.016 U	MG/KG	Chrysene
0.020 U	MG/KG	Benzo(b)Fluoranthene
0.016 U	MG/KG	Benzo(k)Fluoranthene
0.016 U	MG/KG	Benzo-a-Pyrene
0.018 U	MG/KG	Indeno (1,2,3-cd) Pyrene
0.018 U	MG/KG	Dibenzo(a,h)Anthracene
0.016 U	MG/KG	Benzo(ghi)Perylene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

Sample 6101 FY 2006 Project: 06-0436

Extractables Scan

Facility: Gulf States Creosoting

Program: SF

10 U

UG

Phenanthrene

Id/Station: QA-DIB1 / Media: DRY ICE BLANK

Requestor: Donna Webster Project Leader: PMEYER Flowood, MS Beginning: 06/08/2006

Ending:

Produced by: Revell, Dennis

RESULTS UNITS **ANALYTE** RESULTS UNITS **ANALYTE** 10 U UG bis(2-Chloroethyl) Ether 10 U UG Anthracene 10 U UG Benzaldehyde 10 U UG Carbazole 10 U UG Hexachloroethane 10 U UG Di-n-Butylphthalate 10 U UG bis(2-Chloroisopropyl) Ether 10 U UG Fluoranthene 10 U UG n-Nitroso di-n-Propylamine 10 U-UG Pyrene 10 U UG Acetophenone UG Benzyl Butyl Phthalate 10 U 10 U UG Nitrobenzene UG bis(2-Ethylhexyl) Phthalate 10 U 10 U UG Hexachlorobutadiene 10 U UG Benzo(a)Anthracene 10 U υG Caprolactam 10 U UG Chrysene 10 U UG 2-Methylnaphthalene 10 U UG 3.3'-Dichlorobenzidine 10 U UG 1.2.4-Trichlorobenzene 10 U ÚG Di-n-Octylphthalate 10 U υG Naphthalene 10 U UG Benzo(b)Fluoranthene 10 U υG 4-Chloroaniline 10 U UG Benzo(k)Fluoranthene 10 U UG bis(2-Chloroethoxy)Methane 10 U UG Benzo-a-Pyrene 10 U υG Isophorone 10 U UG Indeno (1,2,3-cd) Pyrene 10 U UG Hexachlorocyclopentadiene (HCCP) 10 U UG Dibenzo(a,h)Anthracene 10 U UG 1,1-Biphenyl 10 U UG Benzo(ghi)Pervlene 10 U υG 2-Chloronaphthalene 10 U UG 2-Chlorophenol 10 U υG 2-Nitroaniline 10 U 2-Methylphenol UG 10 U UG Acenaphthylene 10 U IJG (3-and/or 4-)Methylphenol 10 U UG Acenaphthene UG 10 U 2-Nitrophenol 10 U UG Dimethyl Phthalate 10 U UG Phenol 10 U UG Dibenzofuran 10 U UG 2,4-Dimethylphenol 10 U UG 2.4-Dinitrotoluene 10 U υG 2,4-Dichlorophenol 10 U UG 2,6-Dinitrotoluene 10 U UG 2.4.6-Trichlorophenol 10 U UG 3-Nitroaniline 10 U UG 2,4,5-Trichlorophenol 10 U UG 4-Chlorophenyl Phenyl Ether 10 U UG 4-Chloro-3-Methylphenol 10 U UG 4-Nitroaniline 20 U UG 2,4-Dinitrophenol 10 U UG Fluorene 20 U UG 2-Methyl-4,6-Dinitrophenol 10 U UG Diethyl Phthalate 20 U ΠŒ Pentachlorophenol 10 U UG n-Nitrosodiphenylamine/Diphenylamine 20 U UG 4-Nitrophenol 10 U UG Hexachlorobenzene (HCB) 2,3,4,6-Tetrachlorophenol 10 U UG 10 U UG Atrazine 10 U UG 4-Bromophenyl Phenyl Ether

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

6102 FY 2006

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 08/16/2006 14:39

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Project: 06-0436

Program: SF

Sample

Id/Station: QA-DIB2 / Media: DRY ICE BLANK Produced by: Revell, Dennis Requestor: Donna Webster Project Leader: PMEYER Beginning: 06/08/2006

Ending:

RESULTS	UNITS	ANALYTE		·	RESULTS	UNITS	ANALYTE	
10 U	UG	bis(2-Chloroethyl) Ether	•		10 U	UG	Anthracene	
10 U	UG	Benzaldehyde			10 Ŭ	UG	Carbazole	
10 U	UG	Hexachloroethane			10 U	UG	Di-n-Butylphthalate	
10 U	UG.	bis(2-Chloroisopropyl) Ether	•	•	- 10 U	ÜĞ	Fluoranthene	
10 U	UG	n-Nitroso di-n-Propylamine			10 U	UG	Pyrene	•
10 U	UG	Acetophenone			10 Ü	υG	Benzyl Butyl Phthalate	
10 U	UG	Nitrobenzene			10 U	UG	bis(2-Ethylhexyl) Phthalate	
10 U	UG	Hexachlorobutadiene		•	10 U	UG	Benzo(a)Anthracene	
10 U	UG	Caprolactam			10 U	UG	Chrysene	
10 U	UG	2-Methylnaphthalene			10 U	UG	3,3'-Dichlorobenzidine	
10 U	UG	1,2,4-Trichlorobenzene			10 U	UG	Di-n-Octylphthalate	
10 U	UG	Naphthalene			10 U	υG	Benzo(b)Fluoranthene	
10 U	UG	4-Chloroaniline			. 10 U	UG	Benzo(k)Fluoranthene	
10 U	UĠ	bis(2-Chloroethoxy)Methane			10 U	UG	Benzo-a-Pyrene	
10 U	UG	Isophorone			10 U	UG	Indeno (1,2,3-cd) Pyrene	
10 U	UG	Hexachlorocyclopentadiene	(HCCP)		10 U	UG	Dibenzo(a,h)Anthracene	
10 U	UG	1,1-Biphenyl			10 U	UG	Benzo(ghi)Perylene	
10 U	UG	2-Chloronaphthalene			10 U	UG	2-Chlorophenol	
10 U	UG	2-Nitroaniline	•		10 U	UG	2-Methylphenol	
10 U	UG	Acenaphthylene			10 U	UG	(3-and/or 4-)Methylphenol	
10 U	UG	Acenaphthene			. 10 U	UG	2-Nitrophenol	
. 10 U	UG	Dimethyl Phthalate			10 U	UG	Phenol	
10 U	UG	Dibenzofuran			10 U	UG	2,4-Dimethylphenol	
10 U	UĢ	2,4-Dinitrotoluene			. 10 U	UG	2,4-Dichlorophenol	
10 U	UG	2,6-Dinitrotoluene	•		10 U	UG	2,4,6-Trichlorophenol	
10 U	UG	3-Nitroaniline	•		. 10 U	UG	2,4,5-Trichlorophenol	
10 U	UG	4-Chlorophenyl Phenyl Ethe	r <sup>'</sup>		10 U	UG	4-Chloro-3-Methylphenol	
10 U	UG	4-Nitroaniline			20 U	UG	2,4-Dinitrophenol	
10 U	UG -	Fluorene		•	20 U	UG	2-Methyl-4,6-Dinitrophenol	
10 U	ŲG	Diethyl Phthalate	•		20 U	UG	Pentachlorophenol	
. 10 U	UG	n-Nitrosodiphenylamine/Dip	nenylamine	•	20 U	UG	4-Nitrophenol	
10 U	UG	Hexachlorobenzene (HCB)			10 U	UG	2,3,4,6-Tetrachlorophenol	
10 U	UG	Atrazine	•				•	
10 U	UG	4-Bromophenyl Phenyl Ethe	r					•
10 U	UG	Phenanthrene						

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 08/16/2006 14:39

6100 FY 2006 Sample **Extractables Scan** 

Project: 06-0436

Facility: Gulf States Creosoting

Flowood, MS

Requestor: Donna Webster Project Leader: PMEYER Beginning: 06/08/2006 08:45

Ending:

Produced by: Revell, Dennis

Id/Station: QA-BB5 / Media: DRY ICE BLANK

Program: SF

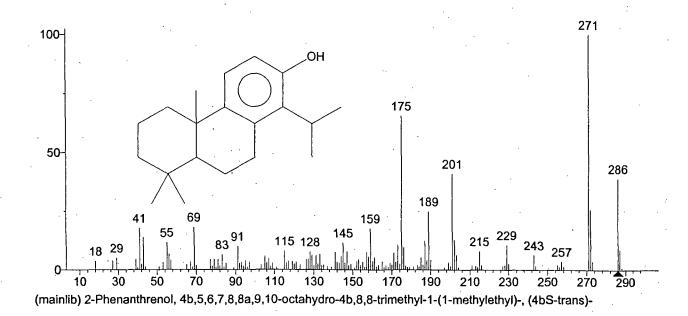
RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
10 U	UG	bis(2-Chloroethyl) Ether	10 U	UG	Anthracene
10 U	UG	Benzaldehyde	10 Ŭ	UG	Carbazole
10 U	UG	Hexachloroethane	10 U	ÚĞ -	Di-n-Butylphthalate
10 U	UG	bis(2-Chloroisopropyl) Ether	10 U	UG	Fluoranthene
10 U	UG	n-Nitroso di-n-Propylamine	10 U	UG	Pyrene
10 U	UG	Acetophenone	10 U	UG	Benzyl Butyl Phthalate
10 U	UG	Nitrobenzene	10 U	UG	bis(2-Ethylhexyl) Phthalate
10 U	UG	Hexachlorobutadiene	10 U	UG	Benzo(a)Anthracene
10 U	UG	Caprolactam	10 U	UG	Chrysene
10 U	UG	2-Methylnaphthalene	10 U	UG	3,3'-Dichlorobenzidine
10 U.	UG	1,2,4-Trichlorobenzene	10 U	UG	Di-n-Octylphthalate
10 U	υG	Naphthalene	10 U	UG	Benzo(b)Fluoranthene
10 U	UG	4-Chloroaniline	10 U	UG	Benzo(k)Fluoranthene
: 10 U	UG	bis(2-Chloroethoxy)Methane	10 U	UG	Benzo-a-Pyrene
10 U	UG	Isophorone	10 U	UG	Indeno (1,2,3-cd) Pyrene
10 U	UG	Hexachlorocyclopentadiene (HCCP)	10 U	UG	Dibenzo(a,h)Anthracene
10 U	UG	1,1-Biphenyl	10 U	ŲG	Benzo(ghi)Perylene
10 U	UG	2-Chloronaphthalene	10 U	UG	2-Chlorophenol
10 Ü	UG	2-Nitroaniline	10 U	UG	2-Methylphenol
10 U	UG	Acenaphthylene	10 U	UG	(3-and/or 4-)Methylphenol
10 U	UG	Acenaphthene	10 U	UG	2-Nitrophenol
10 U	UG	Dimethyl Phthalate	10 U	UG	Phenol
10 U	UG	Dibenzofuran	10 U	UG	2,4-Dimethylphenol
10 U	UG	2,4-Dinitrotoluene	10 U	UG	2,4-Dichlorophenol
10 U	UG	2,6-Dinitrotoluene	10 U	UG	2,4,6-Trichlorophenol
10 U	UG	3-Nitroaniline	10 U	UG	2,4,5-Trichlorophenol
10 U	UG ,	4-Chlorophenyl Phenyl Ether	10 U	UG	4-Chloro-3-Methylphenol
10 U	UG	4-Nitroaniline	20 U	UG	2,4-Dinitrophenol
10 U	UG	Fluorene	20 U	UG	2-Methyl-4,6-Dinitrophenol
10 U	UG	Diethyl Phthalate	20 U	UG	Pentachlorophenol
·10 U	UG	n-Nitrosodiphenylamine/Diphenylamine	20 U	UG	4-Nitrophenol
10 U	UG	Hexachlorobenzene (HCB)	10 U	UG	2,3,4,6-Tetrachlorophenol
10 U	UG	Atrazine		•	
10 U	UG	4-Bromophenyl Phenyl Ether			
10 U	UG	Phenanthrene			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Polytemate is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporting limit. | UJ-Analyte not detected at or above reporti N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.



Name: 2-Phenanthrenol, 4b,5,6,7,8,8a,9,10-octahydro-4b,8,8-trimethyl-1-(1-methylethyl)-, (4bS-trans)-

Formula: C<sub>20</sub>H<sub>30</sub>O

MW: 286 CAS#: 511-15-9 NIST#: 42592 ID#: 96141 DB: mainlib

Other DBs: NIH

Contributor: R RYHAGE MS-LAB KAROLINSKA INSTITUTET STOCKHOLM SWEDEN

10 largest peaks:

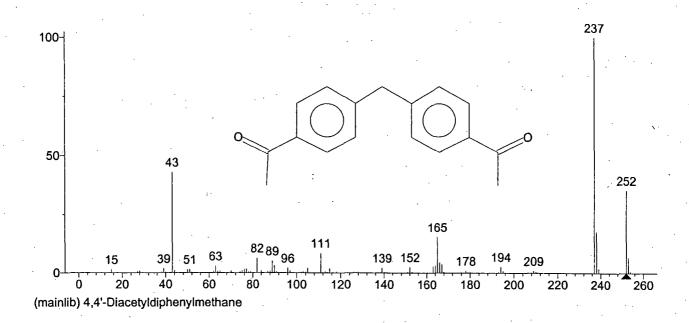
271 999 | 175 653 | 201 409 | 286 391 | 272 258 | 189 249 | 69 182 | 41 178 | 159 177 | 43 140 | Synonyms:

1.Podocarpa-8,11,13-trien-13-ol, 14-isopropyl-

2.Totarol

3.Totarol (b637832k176)

4.(+)-Totarol



Name: 4,4'-Diacetyldiphenylmethane

Formula: C<sub>17</sub>H<sub>16</sub>O<sub>2</sub>

MW: 252 CAS#: 790-82-9 NIST#: 135072 ID#: 90773 DB: mainlib

Other DBs: None

Contributor: NIST Mass Spectrometry Data Center, 1994

10 largest peaks:

237 999 | 43 428 | 252 353 | 238 177 | 165 154 | 111 85 | 253 66 | 82 63 | 89 51 | 166 45 |

Synonyms: no synonyms.

# **APPENDIX II**

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Field Sampling Plan Quality Assurance Plan Data Quality Objectives

## Field Sampling Plan/ Quality Assurance Project Plan

**Gulf States Creosoting Company, Inc.** Flowood, Rankin County, Mississippi

Prepared By

Phyllis Meyer
U.S. EPA Region 4
Science and Ecosystems Support Division
Athens, GA

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To collect the required information, the sampling must be done when the water level conditions are low and there have been no major releases from the Ross Barnett Reservoir to the north of the site. The study will be conducted during the week of May 29, 2006. The sampling date is subject to change based on weather and field conditions. Personnel from the Mississippi Department of Environmental Quality (MDEQ) will assist EPA and the Integrated Laboratory Systems (ILS) with the collection and processing of samples.

## **QUALITY OBJECTIVES/CRITERIA FOR MEASUREMENT DATA**

The data quality objectives for this study were developed using the Guidance for the Data Quality Objectives Process (US EPA 2002) are provided in Appendix A.

## SPECIAL TRAINING REQUIREMENTS/CERTIFICATION

Proficiency with the use of electro-fishing equipment, both backpack shocker and boat shocking units, will be necessary for this field study. Personnel participating in this study will receive an on-site safety briefing prior to initiation of sampling activities. Pertinent safety information is available in Appendix B of this document.

#### DOCUMENTATION AND RECORDS

A bound field logbook will be used for recording information pertinent to this study. All field notes will include information outlined in Section 2.5 of the <u>Ecological Assessment Standard Operating Procedures and Quality Assurance Manual</u>, January 2002 (EASOPQAM). Upon completion of sampling activities, all documents/records obtained during the field investigation will be organized, labeled, and maintained by the project leader during preparation of the report. Upon completion of the report, project records will be submitted to the SESD records room.

The final data report will include data for each media sampled. Tables and maps will also be included in the final report. The text of the report will describe the study collection effort and findings for each station and will include any problems encountered or other noteworthy information. Field data logs will not be included in the final report.

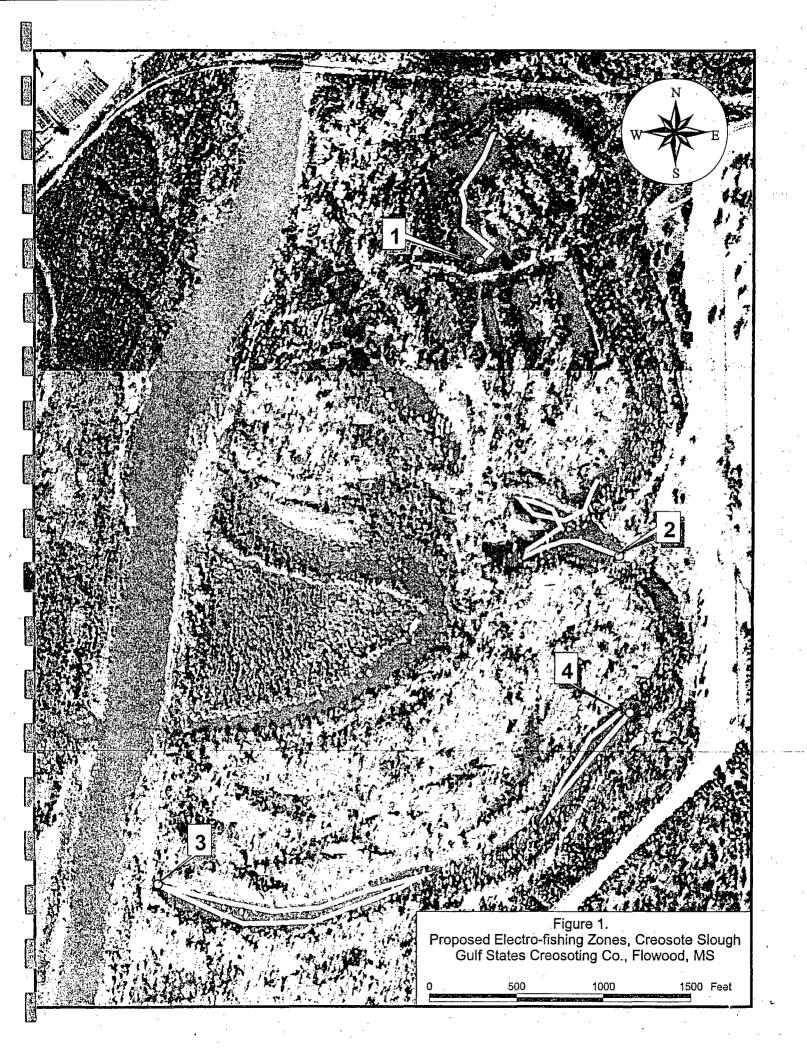
## QUALITY ASSURANCE AND QUALITY CONTROL

Quality control (QC) procedures will be used in the field and laboratory to ensure that reliable data are obtained. During sampling, precautions will be taken to prevent cross-contamination of sampling equipment that could compromise sample integrity. Field and laboratory methods utilized on this project will adhere to USEPA approved guidance and methodology (US EPA, 2005) and the EASOPQAM (US EPA, 2002) and manufacturers instructions. All samples will be handled and custody maintained in accordance with Section 2 of the EASOPQAM. A sonde will be calibrated prior to deployment each day

## REFERENCES

- US EPA. 2005. Analytical Support Branch Laboratory Operations and Quality
  Assurance Manual. US Environmental Protection Agency, Region 4, Science and
  Ecosystem Support Division, Analytical Support Branch, Athens, GA.
- US EPA. 2003. Draft Preliminary Assessment/Site Inspection Report, Gulf States
  Creosoting Company prepared by Weston Solutions, Inc, expressly for EPA.
- US EPA. 2002. Ecological Assessment Standard Operating Procedures and Quality Assurance Manual. US Environmental Protection Agency, Region 4, Science and Ecosystem Support Division, Ecological Assessment Branch, Athens, GA
- US EPA. 1994. Guidance for the Data Quality Objectives Process. EPA QA/G-4. US Environmental Protection Agency, Office of Research and Development, Washington, D.C. EPA/600/R-96/055.

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## DQO FOR FISH, SEDIMENT AND WATER SAMPLING GULF STATES CREOSOTING, FLOWOOD, MS

DQO	STEPS	FISH/SEDIMENT/SURFACE WATER SAMPLING
State the Problem:  * Identify the members of the planning team.  *Identify the primary decision maker of the planning team.		Mississippi Department of Environmental Quality (MDEQ) regulators are concerned that individuals fishing in an area adjacent to the former Gulf States Creosoting Company located at 1625 Flowood Drive (Mississippi Hwy 468), Flowood, Rankin County, Mississippi, may be exposed to elevated levels of PAHs by consuming fish from Creosote Slough.
	*Define problem.  *Specify the available resources and relevant deadlines for the study	Members of the planning team were Donna Webster, U.S. Environmental Protection Agency (EPA), Remedial Project Manager (RPM), Atlanta, GA, Kevin Koporec, EPA Toxicologist, Atlanta, GA, Phyllis Meyer (EPA biologist), field project leader, Athens, GA, Phillip Weathersby, MDEQ, Jackson, MS, Richard Ball, MDEQ.
:		The primary decision maker will be Ralph Howard, EPA (RPM).
		Phyllis Meyer will be responsible for planning and conducting the field investigation; coordinating analytical requirements and compiling the raw data. The field work must be conducted in a one week time frame and is scheduled for May 30 through June 2,
		2006, with a reconnaissance April 10-12, 2006. If this study proceeds as projected and the appropriate data are acquired, analytical results should be available by late August. All project deliverable and task dates are estimates based on analytical laboratory schedules. New information, additional tasks, and changes in scope may result in
		revisions to these dates. The Science and Ecosystem Support Division (SESD) will provide a minimum of two people, a back-pack fish shocking unit, and a boat with electro-fishing capabilities if needed, and other necessary supplies. MDEQ will provide at least one person to assist with the study, lab space to fillet the fish samples
		and an electro-fishing boat if one is available for the study.

2	Identify the Decision *Identify the principal study question. *Define the action that could result from resolution of the principal study question.	Are the body burdens in fish tissue in selected segments of Creosote Slough, adjacent to the former Gulf States Creosoting Company, above the 1E-5 screening value for PAHs (polycyclic aromatic hydrocarbons)? If the data shows that the tissue numbers are above the screening value, then more calculations of the actual risk will be done. If the data shows no exceedance of PAHs, then there is a possibility that no further remedial action will be planned.
3	Identify the Inputs to the Decision  *Identify the information that will be required to resolve the decision statement.  *Determine the sources for each item of information identified.  *Identify the information that is needed to establish the action level.  *Confirm that analytical methods exist to provide the data.	No previous fish data exists for this site.  Fish, sediment and water samples will be collected from approximately 3 stations in Creosote Slough which is located adjacent to the former Gulf States Creosoting facility and one background site located on the northern end of the property, south of the railroad tracks. An attempt will be made to collect edible size fish that are top predators, such as bass or sunfish and fish that are bottom feeders, such as catfish. Depending upon the availability and size of the fish, they will be analyzed either as a composite or as individuals. Composite sediment samples will be collected from each tissue sampling location. The sediment sample will consist of a one-six inch core taken from the beginning, middle and end of each reach that is fished and then composited. One water sample will also be collected from the middle of each fished reach.  Analytical scans for extractable organic compounds (PAHs) will be completed for-all samples. In addition, % lipids and % moisture will be analyzed for all tissue samples. Percent moisture for all sediments.  Lower detection limits (.005 ppm) for all samples will be utilized (this has been negotiated with the chemists and will depend on amounts of tissue available).

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4	Define Study Boundaries *Specify the	1. Determine the concentration of listed contaminants (PAHs) in fish, sediments and water samples.
	characteristics that define the population of interest. *Define the spacial	2. Spacial boundary – Creosote Slough adjacent to the former Gulf States Creosoting, Inc.
	boundary.  *Define the temporal boundary.	3. Temporal boundary – Samples will be collected during the week of May 29 <sup>th</sup> , 2006. Samples will be collected under similar conditions of flow and stage.
	*Define the scale of decision making.	4. Scale of Decision Results from this study will be used to begin making decisions regarding the fishery in Creosote Slough.
	*Identify practical constraints on the data collection	5. Practical Constraints – The study may be delayed if intense rainfall occurs that causes dramatic changes in the flow pattern of the slough due to releases from the Ross Barnett Reservoir.
5	Develop a Decision Rule  * Action Levels of the	Concentrations of PAHs will be used in comparison to background levels.
	study *Develop a decision rule.	If PAHs are detected above background levels, further study may be required to fully characterize the level and specific types of PAHs in the fish tissue.
6	Specify Decision Error Limits *Determine the possible range of the parameters of interest. *Identify the decision errors.	An attempt will be made to collect edible size fish from four different stations. The proposal is to collect a composite of 4-5 top predators and 4-5 bottom feeders from each station. The composites will consist of the same species and will be similarly sized fish. The type and extent of the fishery is not known and deviation from this regime may be altered as the samples are actually collected. All tissues and sediments will be analyzed for PAHs and % moisture. Tissues will also be analyzed for % lipids. Surface water samples will be analyzed for PAHs.

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7	1 ^	timize th sign	e Study		The data will have to be furt possibilit backpack stainless surface.	e a detrin ther eval ty that m c shockir	nental ef uated in ore samp ng and bo	fect on he relation pling ma	uman he to the typy y have to to-fishing	ealth due pes of PA be done g. Sedin	to PAHs AHs dete e. Collections and the contraction in the contraction	s or if the ected. The etion tec Il be coll	e risk fac nere is al hniques ected us	ctors nee lso the will incli ing a	ed lude
			: . :		Sampling 1)	g sites w	ere selec	ted to co	ver the n	navigable	reachs	of Creos	ote Slou	gh (Figu	ire

## APPENDIX III

表

Safety Plan

#### FIELD SAFETY PLAN

1220 3.1	LETTLAN					
SAFETY PLAN						
Site Name: Gulf States Creosoting	Contact: Phillip Weathersby, MDEQ, 601-961-5302					
Address: Motel: Comfort Inn, 235 Pearson Rd	, Pearl, MS					
<b>!</b>	onfirmation #s 198465 and 198466 60.00 Amanda					
Purpose of Visit: Sediment, surface water and fisher	eries work on Creosote Slough, Flowood, MS					
Directions to Site: I-20 West to Hwy 475N or Hwy turn West to Treetops Blvd. for hotel or take Hwy						

#### SITE INVESTIGATION TEAM:

PERSONNEL *	SAFETY CATEGORY	RESPONSIBILITIES
·		
Phyllis Meyer	В	Project leader
Jerry Ackerman	В	ESAT/sampler
Ralph Howard		Regional Project Manager
Richard Ball		MDEQ
Phillip Weathersby		MDEQ
Jerry Banks		MDEQ

All employees have been trained/medically monitored in accordance with OSHA 29 CFR 1910.12 requirements and US-EPA Region IV Field Health and Safety Manual, 1990 edition.

#### PLAN PREPARATION:

Site Safety Officer	Phyllis Meyer	Afreger	Date: 5 26 06
Branch Safety Officer	Phyllis Meyer	Apoly	Date: 5 26 66
Section Chief:	Bobb	, L Carta	Date 5/26/06

#### SITE HAZARDS:

Electro-fishing, Insects, snakes								
					:			

TRAILER LICENSE:

WHERE PARKED: ConSteel

Complete this plan, before going boating and leave it with a reliable person who can be depended upon to notify the Coast Guard, or other rescue organization, should you not return as scheduled. Do not file this plan with the Coast Guard.

PROJECT DATES - 05/30/06 - 06/02/06

(if overnight, date returning) -

1. NAME OF PERSON REPORTING: Phyllis Meyer TELEPHONE NUMBER: 706-338-2867

BOAT MAKE	COLOR	LENGTH	ENGINES	OCCUPANTS
Pontoon	Silver	28	ОВ	
Parker	White	25	2 OB	
Privateer	White	24	2 OB	
Parker	White	23	2 OB	
Parker	White	21	ОВ	
Mako	White	20	ОВ	
Privateer	White	18	ОВ	
Shocker-Schaffer	Tan	18	ОВ	
Boston Whaler	White	17	ОВ	
Shocker-Fisher	Green	16 .	ОВ	
Canoe Old Town X 2	Green	16		
Canoe	Aluminum	16		
Shocker-Scandy/White	Aluminum	15	ОВ	Phyllis Meyer, Jerry Ackerman, Richard Ball (MDEQ)
Jon	Gray	14	ОВ	
Jon X 2	Gray	12	ОВ	
RiverHawk - Canoe	Green	12		

Jon	Gray	14	ОВ		•
Jon X 2	Gray	12	ОВ		
RiverHawk - Canoe	Green	12			
4. TRIP EXPECTATION FROM - Creosote Slow		T - 0800	(TIME)		
GOING TO -Sites 1,2	4,3 in that ord	ler			
EXPECTED TO RET	URN BY: 210	0	(TIME)		<u>.</u>
AND IN NO EVENT  5. IF NOT RETURNED NAME: TELEPHONE NUMBI	BY		(TIME) CALL THE COAS	ST GUARD, OF	R (LOCAL AUTHORITY)
6. SURVIVAL EQUIPM x_PFDs CLOTHING WATER EPIRB	ENT: (CHEC: fla fla		_	MIRROR FOOD xANCHOR	SMOKE SIGNALSx_PADDLESRAFT OR DINGHY
7. RADIO:YES TYPE: 8. ANY OTHER PERTIN	FRI	EQS			

COLOR/MAKE OF AUTO:

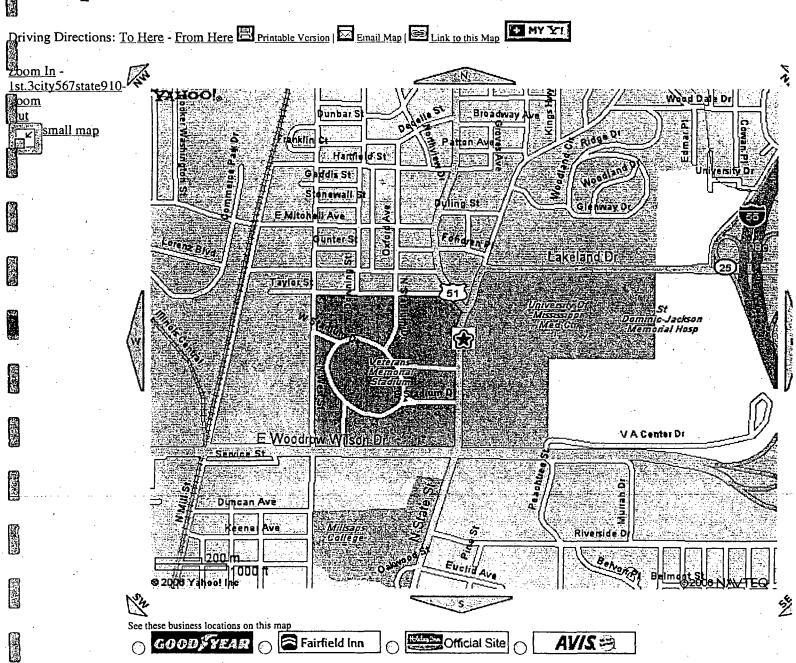
Yahool My Yahool Mail Make Yahool your home page

YAHOO LOCAL Sign In New User? Sign Up

Your trip to the Southeast starts here



## Map for: 2500 N State St Jackson, MS 39216-4500 Sav



Zoom in & Re-Center Re-Center only View Map Legend

SPONSORED LINKS

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Reference 6 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423



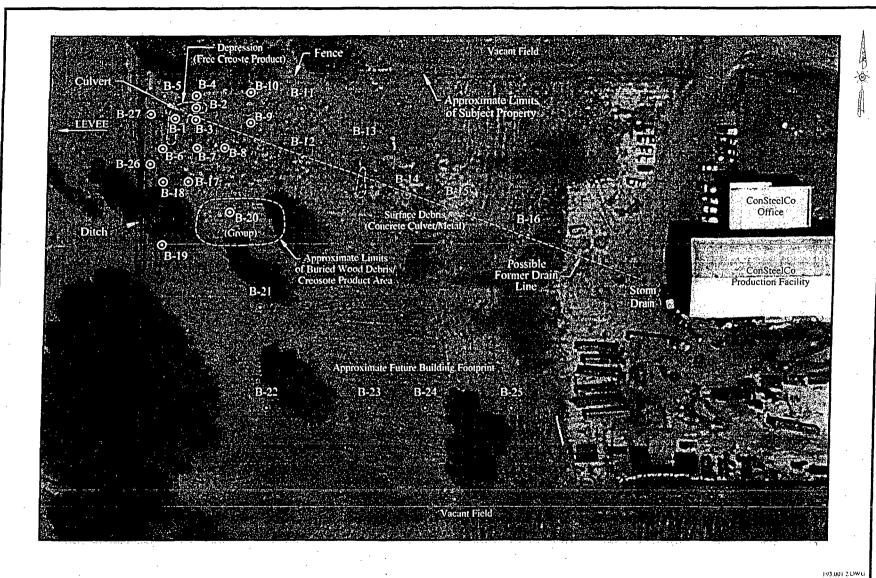
LIMITED SOIL ASSESSMENT CONSTEELCO, INC. 1625 FLOWOOD DRIVE FLOWOOD, MISSISSIPPI

#### PREPARED FOR

MS. KELLY BLACKWOOD
PHELPS DUNBAR
111 EAST CAPITOL STREET, #600
JACKSON, MISSISSIPPI 39201

EarthCon Project Number C193.001

Submitted October 9, 2006



LEGEND:

Soil Boring Location



Soil Boring Location With Creosote Odor

Area of Concern



#### Earth Consulting Group, Inc.

P.O. Box 1246 Madison, Ms 39130 Tel: (601) 853-2134 Fax: (601) 856-3978

- DRAWN BY:		CHECKED BY:	SCALE:	DATE	:	PROJECT NO:
	 Glen Ivey	Jay Ferris	NTS		09/22/06	C193.001
PROJECT:	:	ConSteelCo Inc - 1625	Flowood Drive - Fl	owoo.	1 Mississint	ni

SOIL BORING LOCATION PLAN

#### **EXECUTIVE SUMMARY**

Earth Consulting Group, Inc. (EarthCon) was authorized by Ms. Kelly Blackwood of Phelps Dunbar to perform a limited soil assessment of the western portion of the ConSteelCo property located at 1625 Flowood Drive, Flowood, Rankin County, Mississippi. The limited soil assessment included installation of soil borings and selected soil sampling and analysis to assess the area for the presence of creosote compounds related to former Gulf State Creosoting, Inc. site operations that occurred prior to purchase of the property by ConSteelCo.

Twenty-seven (27) soil borings were installed to a depth of approximately 8.0 feet below surface grade in the western portion of the subject property. The borings were visually described and soil conditions were recorded on boring logs. Borings with obvious visual or olfactory indications of creosote contamination were noted, and an approximate delineation of the lateral extent of the residual creosote impacts was achieved. Approximately 15,000 square feet of the northwest corner of the property appeared to be impacted with creosote odors and/or soil staining in the shallow subsurface.

Free creosote product was observed in soil borings at two (2) locations on the site. These areas included a depression near the northwest corner of the subject property, and an area containing buried wood debris in the west-central portion of the subject property. Borings installed within the footprint of a proposed building addition on the property, and along the route of a possible drain line on site, did not contain evidence of creosote impacts in the shallow soil.

Laboratory analytical results of soil samples collected from each of the two (2) identified creosote locations on site contained concentrations of semi-volatile organic compounds in excess of the Mississippi Department of Environmental Quality (MDEQ) Target Remediation Goals (TRGs) for "Restricted Use" scenarios. Several other semi-volatile organic compounds were detected at concentrations within allowable TRG limits.

The soil conditions identified in this study appear directly related to past site operations by Gulf States Creosoting, Inc. This site has been the focus of preliminary environmental investigations by the U.S. Environmental Protection Agency and the MDEQ. EarthCon recommends consulting with legal counsel and the appropriate environmental regulatory agencies to determine an appropriate course of action for the property.

#### Certification:

Limited Soil Assessment ConSteelCo, Inc. 1625 Flowood Drive Flowood, Mississippi

Prepared for

Mr. Randy Kenner ConSteelCo, Inc. 1625 Flowood Drive Flowood, Mississippi

Earth Consulting Group, Inc. hereby certifies the aforementioned report constitutes an accurate presentation of the investigation, research, and findings developed during the completion of the Limited Soil Assessment prepared for, and submitted to, the client as their approved Engineer of Record.

Signed:

Senior Project Manager

Mississippi Registered Professional Geologist #0045

Michael J. Brady, P.E.

Senior Engineer

Mississippi Registered Professional Engineer #12402

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#### 1.0 INTRODUCTION

Earth Consulting Group, Inc. (EarthCon) was authorized by Ms. Kelly Blackwood of Phelps Dunbar to perform a limited soil assessment of the western portion of the ConSteelCo, Inc. (ConSteelCo) property located at 1625 Flowood Drive in Flowood, Rankin County, Mississippi (see Figure 1 - Site Location Map). The subject property exists within the boundary of the former Gulf States Creosoting, Inc. (Gulf States Creosoting) facility. The Gulf States Creosoting facility has been the object of preliminary environmental investigation by the U.S. Environmental Protection Agency (EPA) and the Mississippi Department of Environmental Quality (MDEQ), including investigation to evaluate inclusion of the property on the National Priorities List for environmental remediation (a.k.a. Superfund). The limited soil assessment in this study was performed to evaluate site conditions prior to a pending commercial transaction for the subject property.

The limited soil assessment included installation of soil borings in the western portion of the property to determine if residual creosote impacts were present, and to laterally delineate any detected impacts to the subsurface. Specifically, the scope of work included:

- Reviewing selected documents outlining the general nature of the property with regard to potential contaminants;
- Locating and marking of underground utilities in the area through the Mississippi One Call service and local utility providers;
- Installing soil borings within the vicinity of a depression near the northwest corner of the subject property;
- Installing soil borings along the route of a potential drain line extending from the northwest corner of the property to a storm grate on the west side of the existing production facility;
- Installing soil borings within the footprint of a proposed building addition on the southwest portion of the subject property;
- Collecting three (3) soil samples from creosote-impacted areas on site to characterize the degree of contamination in areas deemed to have the greatest evidence of creosote;
- Preparing a report documenting the findings of the soil assessment.

#### 2.0 SOIL BORING AND MONITORING WELL INSTALLATION

A limited program of soil boring installation was performed to determine whether creosote had been released to the subsurface during past Gulf States Creosoting operations on the subject property. The boring locations selected and sampling methodology are described in the following sections and boring locations are shown in Figure 2 – 2003 Aerial Photograph - Soil Boring Location Plan.

#### 2.1 Sampling Location Rationale

The western portion of the subject property was observed as a vacant field and appeared regularly maintained (mowed) at the time of the limited soil assessment. This portion of the property has reportedly not been used in site operations since its purchase by ConSteelCo in 1994. The approximately 8-acre subject property was formerly part of the approximately 125-acre Gulf States Creosoting facility which operated several years prior to the purchase by ConSteelCo. Environmental investigations have been conducted in the past on the subject property and adjacent properties by the EPA, MDEQ, and other environmental consulting firms. A recent site investigation related to a pending commercial transaction for the ConSteelCo site discovered a small depression in the northwest corner of the property that contained remnants of a concrete culvert. Soil borings performed by others reportedly detected free creosote product in the shallow subsurface in this area. Soil borings were initially installed by EarthCon in this area to confirm the presence of creosote in the subsurface. Borings B-1 through B-12, B-17 through B-20, B-26, and B-27 were installed in a general radial pattern away from this area to attempt to delineate the lateral extent of the creosote impacts associated with this feature.

Other areas of potential concern on the property included a possible buried drain line leading from the west side of the existing ConSteelCo production building to the concrete culvert remnant in the northwest corner of the property. Soil borings B-13 through B-16 were installed along the suspected route of this potential drain line including areas of stressed vegetation observed along the route.

Soil borings B-21 through B-25 were installed within the footprint of a planned new ConSteelCo production and office facility on the southwest portion of the property. These borings were

installed to assess possible residual creosote impacts to the subsurface beneath the area where workers would be present during planned future operations.

#### 2.2 Sampling Methodology

The assessment objective was to attempt to laterally delineate creosote compounds previously reported in the shallow soil on site. Soil borings were installed to a depth of approximately 8.0 feet below surface grade at 27 locations on the western portion of the property. Due to the inherently distinct odor and dark brown to black color of creosote, odors (olfactory evidence) and visual indications of creosote were utilized to determine whether the soil samples retrieved from a boring location contained creosote compounds. Upon completion of the sampling and visual classification of the 27 soil borings, EarthCon returned to three (3) boring locations (B-2, B-7, and B-20), and installed a boring adjacent to each original boring location to collect a representative sample of the creosote impacted material in these areas. These locations were selected to enable comparison of chemical concentrations of samples deemed as the most impacted soils in the assessment to the MDEQ Tier 1 Target Remediation Goals (TRGs) for "Restricted Use" (industrial) property.

A truck-mounted Model 540 UD Geoprobe<sup>®</sup> direct-push drilling rig (Geoprobe) was used to collect the soil samples in this assessment. The Geoprobe was equipped with a 48-inch long by 2-inch diameter Macro-core sampling device which utilizes acetate inner sleeves to retrieve continuous soil samples as the sampler is pushed through the soil. This process was repeated to yield a continuous soil core to the depth of boring termination. Upon retrieval from the Macro-core sampler, the soil-filled acetate sleeve was cut open and the soil was described and classified on soil boring logs. Noticeable creosote odors and/or staining (if present) were recorded for each soil sampling interval. The boring logs prepared in this limited soil assessment are presented in Appendix A.

Groundwater was not encountered in the soil borings installed in this assessment. No groundwater monitoring wells were installed as part of this limited soil assessment. The soil samples collected in this assessment were generally described as moist, with moisture content noted to increase slightly with depth. A wet sand stratum was encountered at a depth of

approximately 7.5-8.0 feet below surface grade in borings installed in the southeastern portion of the study area in the vicinity of boring B-25. Based on the visual evidence obtained in the soil borings installed in this assessment, EarthCon estimates that the groundwater table exists within 10-20 feet of the surface.

#### 2.3 Laboratory Analysis

Based on the chemical nature of the suspected site contaminants, the three (3) soil samples collected for laboratory analysis in this assessment were analyzed for semi-volatile organic compounds by EPA Method 8270C. Environmental Science Corporation in Mt. Juliet, Tennessee performed the laboratory analysis. The laboratory analysis results were compared to the MDEQ Tier 1 Target Remediation Goals for soil in the "Restricted Use" category as the subject property is an industrial site and is planned for continued industrial use in the future. The Limited Soil Assessment Results are discussed in the following section.

#### 3.0 LIMITED SOIL ASSESSMENT RESULTS

The following sections describe the results obtained in the limited soil assessment for the subject property.

#### 3.1 Soil Stratigraphy

The shallow soil in the western portion of the subject property can be generally described as consisting of a soft to firm, low-plasticity clay from the surface to a depth of approximately 4.0 feet below surface grade. A dense, highly plastic clay layer was generally observed from approximately 4.0 - 6.5 feet below surface grade. A light brown, fine sand stratum was observed from approximately 6.5 - 8.0 feet below surface grade in some portions of the site.

#### 3.2 Creosote Impacted Area

Based on visual and olfactory evidence, creosote impacted soils appear to exist within an area measuring approximately 100 feet (east to west) by approximately 150 feet (north to south) (approximately 15,000 square feet) in the northwest corner of the ConSteelCo property. Additional information concerning creosote-impacted areas on site is presented in the following sections.

#### 3.2.1 Northwest Corner Depression

Free creosote product was observed in two (2) areas on the ConSteelCo property in this assessment. Borings installed in a shallow depression in the northwest corner of the subject property were observed to contain degraded, soft soils and liquid creosote product in the upper few feet, with heavy creosote odors and stained soil to the termination depth of the borings at approximately 8.0 feet below surface grade. This depression appears as an irregular U-shaped area measuring approximately 50 feet long by approximately 15 feet wide (approximately 750 square feet). Noticeable odors and/or creosote staining were observed in several borings surrounding this depression, generally encompassing the northwest corner of the property.

#### 3.2.2 Buried Debris Area

Within the creosote impacted portion of the ConSteelCo property described above, an approximately 75 foot (east to west) by 50 foot (north to south) (3,750 square feet) area

surrounding boring location B-20 was observed to contain buried creosote-soaked wood debris. This debris was encountered in several boring locations at a depth of approximately 3.5 - 4.5 feet below surface grade. The debris was impenetrable using the soil sampling device, although small samples of creosote-soaked wood were collected on occasion from the drive shoe at the bottom of the soil sampler. The lateral extent of this location was assessed by attempting to penetrate the debris in several shallow borings after initially encountering the debris at the boring B-20 location. Once the Geoprobe moved laterally in each cardinal direction enough to avoid the debris area, the soils encountered were similar to native materials observed on site, with slight odors and/or soil staining. The thickness of the buried wood debris in this area was not defined.

#### 3.3 Possible Drain Line Location

A storm drain grate was observed approximately 50 feet west of the western edge of the ConSteelCo production facility building. This drain was reportedly discovered after the purchase of the property by ConSteelCo. The route and extent of the stormwater drainage pipe from this location is unknown, but the general direction of the pipe exiting the drain appears toward the northwest corner of the subject property where a broken concrete culvert remnant was located in the previously mentioned shallow depression. Patches of barren ground indicating stressed vegetation were observed along a line from the drain to the remnant culvert. Borings B-12 through B-16 were installed along a line between the storm drain and the remnant culvert to evaluate the probable pipeline route, should it exist, and assess the areas of stressed vegetation for indications of creosote material.

Boring results from the locations tested in this area did not reveal evidence of the drain line existence or the presence of creosote in the soil samples collected. The route of the storm water drain and the cause of the stressed vegetation remain unknown at this time.

#### 3.4 Future Building Footprint

The ConSteelCo facility reportedly intends to construct an additional T-shaped office and production facility building on the southwestern portion of the property (see Figure 2). Borings B-21 through B-25 were installed adjacent to, or within, the general footprint of the proposed

building location to assess the potential for creosote-impacted soils existing beneath the future building foundation. The soil borings within the proposed construction footprint did not contain visual or olfactory evidence of creosote to a depth of approximately 8.0 feet below surface grade. Soil materials generally consisted of low-plasticity silty clay grading to a highly plastic, dense clay with depth. A layer of fine sand was evident in the bottom few feet of borings B-24 and B-25, increasing in thickness to the east.

#### 3.5 Laboratory Analysis Results

Laboratory analysis of soil samples submitted in this assessment detected creosote-related compounds in each of the areas tested. The chemical constituent naphthalene was detected in each of the samples at concentrations exceeding the MDEQ Target Remediation Goal for soil in Restricted Use settings, with the highest concentration of 4,600 mg/kg exhibited in the sample collected from the area of boring B-20. The soil analytical data is summarized in Table 1. A copy of the laboratory analytical data report is included as Appendix B.

#### 4.0 CONCLUSIONS AND RECOMMENDATIONS

EarthCon has completed a limited soil assessment on the western portion of the ConSteelCo, Inc. property located at 1625 Flowood Drive, Flowood, Rankin County, Mississippi. The results of assessment revealed the northwest corner of the property to be impacted with creosote compounds in the shallow soil, with two (2) locations exhibiting free creosote product in the subsurface. The area of creosote impacts appears to encompass an area of approximately 15,000 square feet to a depth of at least 8.0 feet below ground surface in most areas. Faint residual creosote odors were noted in borings located along the north and western property boundaries; however, the extent of residual creosote impacts has not been fully defined off site.

Other areas of interest in the western portion of the subject property were assessed for indications of creosote impacted soil. These areas included the presumed route of a potential drain line and within the footprint of a proposed building addition on the subject property. Soil samples collected in these areas did not contain visual or olfactory evidence of creosote impacts. Proposed building plans do not appear to be detrimentally impacted by the results of this study with regard to creosote impacted soil. However, potential hazard communication per 29 CFR 1910.1200 will be required based on the results of this study.

The creosote impacted soil on the subject property appears directly related to the former Gulf States Creosoting, Inc. operations. The Gulf States Creosoting, Inc. facility has been the focus of preliminary studies by the EPA and MDEQ. Based on information contained in the supporting documents reviewed by EarthCon, the results of these studies indicated the Gulf States Creosoting, Inc. operations did not meet the minimum scoring criteria to be included on the National Priorities List (Superfund), and a status of "No Further Remedial Action Planned" has been proposed for the site by EPA. This recommendation has not been accepted as yet by the MDEQ until further testing and analysis is conducted on adjacent properties (Creosote Slough).

The results of this assessment indicate an area of buried creosote and related waste material exists in the northwest portion of the ConSteelCo, Inc. property. This waste material appears directly related to former Gulf State Creosoting operations on the property. Soil samples

collected in this assessment contained concentrations of chemical contaminants in excess of MDEQ Target Remediaton Goals for Restricted Use (industrial) property settings.

Based on the information contained in this study, EarthCon recommends ConSteelCo, Inc. consult with legal counsel and appropriate regulatory agencies to determine and document an appropriate future course of action for the facility.

Table 1 Soil Analytical Data Summary

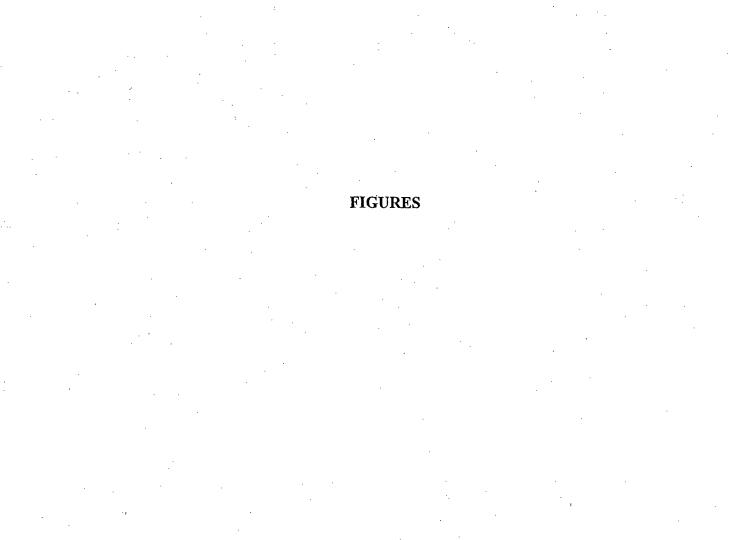
# Table 1 Soil Analytical Data Summary Limited Soil Assessment ConSteelCo, Inc. 1625 Flowood Drive Flowood, Mississippi EarthCon Project No. C193.001

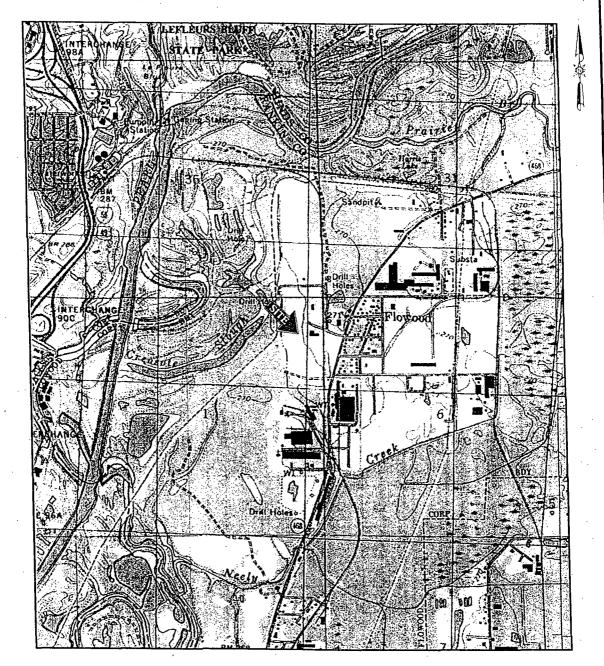
	SOI	L SAMPLES - September 20,		
Sample No.	Location	Detected Compound	Concentration (mg/kg)	MDEQ TRG* (mg/kg)
B-2 (5.5 feet)	Northwest Corner	Acenaphthene	81	123,000
<u> </u>		Fluoranthene	120	81,700
		Fluorene	75	81,700
		Naphthalene	770	247
		Phenanthrene	200	61,300
	· · ·	Pyrene	100	61,300
B-7 (8.0 feet)	Northwest Corner	Acenaphthene	390	123,000
	· · · · · · · · · · · · · · · · · · ·	Anthracene	120	613,000
		Benzo (a) anthracene	82	7.8
•	4 · · · · · · · · · · · · · · · · · · ·	Fluoranthene	490	81,70
•	•	Fluorene	370	81,70
		Naphthalene	2,300	24
		Phenanthrene	1,000	
		Pyrene	330	61,30
B-20 (5.0 feet)	Buried Debris Area	Acenaphthene	2,400	123,00
		Anthracene	5,500	613,00
	:	Benzo (a) anthracene	930	7.8
1		Chrysene	850	78
		Fluoranthene	5,400	81,70
· ·		Fluorene	3,300	81,7
		Naphthalene /	4,60	
		Phenanthrene	7,40	
		Pyrene	3,20	

Notes.

MDEQ TRG\* - Mississippi Department of Environmental Quality Target Remediation Goal for "Restricted Use" (industrial) Soil Setting

Compounds Highlighted in Yellow Exceed the Established MDEQ TRG







SOURCE: USGS 7.5' MAP - JACKSON QUADRANGLE - 1998

Earth Consulting Group, Inc.



	<u> </u>
	•
P.O. Box 1246 Madison, MS 39130	Tel:(601)853-2134 Fax:(601)856-3978

Glen Ivey Jay Ferris 1" = 2000' 09/21/06 C193.001

PROJECT: ConSteelCo, Inc. - 1625 Flowood Drive - Flowood, Mississippi

SITE LOCATION MAP

FIGURE

Figure 2 2003 Aerial Photograph - Soil Boring Location Plan







LEGEND:

Soil Boring Location



Soil Boring Location With Creosote Odor



Area of Concern

B-20 (Group) - Area of Numerous Boring Attempts to Penetrate Buried Wood Debris

#### Earth Consulting Group, Inc.

P.O. Box 1246 Madison, Ms 39130 Tel: (601) 853-2134 Fax: (601) 856-3978

DRAWN BY: Glen Ivey C193 001 PROJECT: ConSteelCo, Inc - 1625 Flowood Drive - Flowood, Mississippi FIGURE.

2003 AERIAL PHOTOGRAPH - SOIL BORING LOCATION MAP



Appendix A Soil Boring Logs

Projec	t N	0.	C19	3.001	Facility ID	ConSteelCo	Date Drilled	_	9/19/06	To	tal Depth	8.0 ft.
Site 1625 Flowood Drive, Flowood, Mississippi						Start Time	_	0900	Time C	ompleted	0915	
Projec	Project Name Limited Soil Investigation						Drill Rig	_				
Geolo	Geologist Jay Ferris Driller D. Riley								Macı	oCore 48-	inch Direc	t Push
Boreb	ole	Co	mpletion	/Abando	onment	Borehole allowe	d to collapse to s	urfa	ice grade			
eet)	ater	ype			<del></del>							
Depth (feet)	Groundwater	Sample Type	USCS Symbol		Mat	erial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
	Demons											
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\ <u>\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\</u>	1							N	1	-		<del> </del>
	1_						•	R	I.			<u> </u>
	l										`	
5.0	-								<b></b>			<del> </del>
1							·					
	T					<del></del>		_{//				
6.0	-		СМ	Greenis	sh brown, firm,	silty CLAY (moist	) (heavy creosote		<b></b>			<u> </u>
				Jugary						!	· ·	1
	$\dagger$			<del> </del>				{/	<b></b>	<del></del>		
7.0		_	W. Fried						<b>//</b>			
			SP		rown, fine SAN te staining)	D (moist) (heavy c	reosote odor) (hea	vy				
+	+				co scanning)				<b>//</b>			<del> </del>
8.	0				· · · · · · · · · · · · · · · · · · ·							
					Boring 8.0 feet	below surface grade						
	82		ample Ty		_	Groundwat					mments	
	_			ous Core	<b>!</b>	Detected Groundw			*:		r laboratory	analysis
	_		Split Sp		E	quilibrated Water Lo	evel V		ppm:	parts per m		
			Hand A	uger					NR:	no recover	у	•

Proje	ct N	ło.	C19	3.001	Facility ID	ConSteelCo	Date Drilled		9/19/06	To	tal Depth_	8.0 ft.		
Site	ite 1625 Flowood Drive, Flowood, Mississippi						Start Time		0915	Time C	ompleted_	0925		
Proje	Project Name Limited Soil Investigation						Drill Rig	Geoprobe 540 UD						
Geole	ogis	t _	Jay Fo	erris	Driller	D. Riley	Auger Type	pe MacroCore 48-inch Direct Push						
Bore	hole	e Co	mpletion	/Abandon	ıment	Borehole allowe	d to collapse to s	urfa	ice grade					
eet)	vater	Type												
Depth (feet)	Groundwater	Sample 7	USCS Symbol	<del></del>	Mate	erial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time		
 					:						·			
								N						
1.0			CL	Greenish	brown, silty, p	olastic CLAY (moi	st) (creosote odor)	R						
					÷	••					<del>-</del>			
2.0	-				· · · · · · · · · · · · · · · · · · ·		<u> </u>							
<u></u>	_		Debris	   Debris P	vooden (railro	ad tie?) (creosote ¡	reduct and adar)							
3.0			Debris	Debi is,	· · · · · · · · · · · · · · · · · · ·	ad de.) (ereosote	or oduct and odor)							
3.0	╁			<u> </u>		·		W		<del></del>	<del> </del>	<u> </u>		
	1		sc	Dark brown, slightly clayey, fine SAND (moist) (slight to						<b></b>				
4.0				moderate	oderate creosote odor)									
	T				······································			N	1			<del> </del>		
-	+							R	<u></u>		<del> </del>	·		
5.0	)		CII	Dank		,	• • • • • • •				Ì			
			CH	Dark gra	y, still, dignly	plastic CLAY (mo	ist) (creosote odor	) <i>[[</i>				10:30		
	+	-						<b></b>	*B-2	8270C	(9/20/06)			
6.0	0			·			· 			·				
7.	١	i i		Reddish	hrown élight	ly candy friable l	ow-plasticity CLAY				ŀ			
	+		CL		eavy creosote		m-plasticity CLA		<b></b>			<del>-  </del>		
									<b>//</b>		<u> </u>			
8	.0	20.40		End of h	Coring Off Cade	palou surface	<u> </u>							
十			Sample Ty		ornig a.v icet	pelow surface grade Groundwat		十		Co	mments			
	Į			uous Core	First	Detected Groundw	<del></del> .		*:		laboratory	analysis		
			Split Sp	poon	E	quilibrated Water L	evel V		ppm:	parts per m				
	Ī	Ш	Hand A	Auger			•	-	NR:	no recover	у			
L					<u> </u>	·	· · · · · · · · · · · · · · · · · · ·		<del> </del>	<del> </del>				

Proj	ect l	No.	C19	3.001	Facility ID	ConSteelCo	Date Drilled	_	9/19/06	To	tal Depth_	8.0 ft.		
Site		1	625 Flov	vood Drive	e, Flowood, N	Start Time		0945	Time C	ompleted _	1000			
Pro	Project Name Limited Soil Investigation								Geoprobe 540 UD					
Geo	Geologist Jay Ferris Driller D. Riley Auger Type								Macr	oCore 48-	inch Direc	Push		
Bor	Borehole Completion/Abandonment Borehole allowed to collapse to surf													
(eet)	Groundwater	Type			,									
Depth (feet)	pun	Sample,	USCS					Recovery	PID Reading	Sample	Analytical	Sampling		
Dep	S.	Sam	Symbol		Mate	erial Description		Reco	(ppm)	Collected	Method	Time		
						,		N						
	+			Daule busin	aliabeta as	.a., 6.5.11. 1	11.14 CK 137	R				ļ		
1.0			CL	(moist) (ne		idy, friable, low-pla	isticity CLA I				1			
<u> </u>	1				,				<del></del>	<del> </del>				
<u> </u>	$\perp$													
			Debris		thtly cemented	l sand (debris) (ligh	t creosote odor)		·					
2.		- 8		(dry)	<del></del>	· <u></u>	<u> </u>			<del> </del>		<del> </del>		
1				ļ						1		1		
$\vdash$	+				;					<del> </del>	1			
3.	0					•								
L											<u> </u>			
												]		
4	.0			1			•					· ·		
				ŀ	•									
$\vdash$	+	-		Gray to	reddish brown	, mottled, stiff, fria	hle low-plasticity				<del></del>	+		
5	.0		CL		lry) (slight odd		bic, to w-plasticity				1			
	1			į						1	1			
L		_				•			<b></b>					
Ľ	5.0				•						<u> </u>	<del> </del>		
				-						1	c			
-	-						•		<b>3</b>	<del></del>		<del></del>		
1	7.0													
				ļ								1		
				<u> </u>	·····	·			<b>2</b>					
	8.0		СН		n brown to gra creosote odor)	y, mottled, stiff, pla	stic CLAY (moist							
ŀ	0.0	<u>!</u>	<b>98</b> ()			below surface grade		-14	<u>//</u>					
t			Sample T			Groundwate	<u></u>	$\top$		<u>Co</u>	mments	<u>-</u>		
			Contin	uous Core	First	t Detected Groundwa	ater V		*:	selected for	r laboratory	analysis		
			Split S	poon	1	quilibrated Water Le		1	ppm:	parts per m		-		
1					_			1	NR:	no recover				
	Hand Auger								Tite no receivery					

roject	No.	C193	3.001	Facility ID	ConSteelCo	Date Drilled	9/19/06	_ To	tal Depth _	8.0 ft.		
ite 1625 Flowood Drive, Flowood, Mississippi				Start Time	1000	Time C	Time Completed _					
roject Name Limited Soil Investigation					Drill Rig		Geoprobe 540 UD					
Geolog	ist	Jay Fe	rris	Driller	D. Riley	Auger Type	Mac	roCore 48-	inch Direc	t Push		
Boreho	ole Co	mpletion/	'Abando	nment	Borehole allowe	d to collapse to s	urface grade	<u> </u>	·	···		
eet)	Type			<del></del>	<del></del>							
Depth (feet)	Sample 7	USCS Symbol		Mate	erial Description		Reading (ppm)	Sample Collected	Analytical Method	Sampling Time		
		.		;								
								_				
1.0	-						<b></b>	<u> </u>	<del>                                     </del>			
			•			•						
								<b></b>				
2.0	_			:			<b>//</b>	<del></del>	<del> </del>	<del> </del>		
				ς								
			•					<del></del>	<del>                                     </del>	<del> </del>		
3.0	_							· · · · · · · · · · · · · · · · · · ·				
			•		:				1	1.		
┟╼╼┼	-	СН	Gray, de	nse, plastic CL	AY (dry to moist)	(no odor)	<b>//</b>		<del> </del>			
4.0				:		•			1			
	_	200	Ì	٠.		•	<b>//</b>					
5.0												
							<b>//</b>	-				
								1				
6.0					•		<b>%</b>			<del></del>		
			ļ					}				
7.0			<b>_</b>		<u>-                                      </u>					_		
1			Light a	rov firm plact	ic CLAY (moist) (l	iaht ta madayata						
-		СН	creosote		ic CLAT (moist) (i	igut to moderate	<b>//</b>		-	<del>-                                     </del>		
8.0	65											
-		Sample To		Boring 8.0 feet b	elow surface grade							
1		Sample Typ Continue		Ei	Groundwat		_		mments	onolyais		
					Detected Groundwa quilibrated Water Le		1 .	*: selected for laboratory analysis				
		Split Spo		E	jumbrated water Le	VCI V	ppm:					
1		Hand A	nger				NR:	no recover	у			

Project	No.	C19	3.001 F	acility ID	ConSteelCo	Date Drilled	_9	/19/06	To	tal Depth_	8.0 ft.		
Site	1	625 Flow	ood Drive,	Flowood, M	ississippi	Start Time		1015 Time Completed 1025					
Project	t Nam	e	Limited	d Soil Investi	gation	Drill Rig	_	Geoprobe 540 UD					
Geolog	gist _	Jay Fe	rris	Driller	Auger Type		Macr	oCore 48-	inch Direc	Push			
Boreho	ole Co	mpletion	/Abandon	ment	Borehole allowed	l to collapse to s	urfac	e grade					
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time		
1.0				: .				···					
			Light gray (CLAY (dry		wn, mottled, dry,	slightly plastic		<del></del>					
2.0			l	:									
3.0										·			
				•		•							
4.0											<u> </u>		
5.0													
5.0		СН	Gray to re (no odor)	ddish brown,	mottled, dense, pl	astic CLAY (dry)				-	<del> </del>		
6.0								·					
											-		
7.0								<del></del>					
8.0		SM/SF	Light gra		yey fine SAND to f	ine SAND (moist)							
	1		End of Bo		elow surface grade								
-	S	Sample Ty	pe		Groundwate	r			Co	mments			
		Continu	ous Core	First	Detected Groundwa	ter ▼		*:	selected for	· laboratory	nalysis		
		Split Sp Hand A		Eq	uilibrated Water Le	vel ∇		ppm: NR:	parts per m				

Project N	lo. <u>C1</u>	93.001 F	acility ID	ConSteelCo	Date Drilled	9	/19/06	Tot	al Depth_	8.0 ft.
Site	1625 Flox	wood Drive,	, Flowood, N	<b>Sississippi</b>	Start Time	. <del>-</del>	1030	Time C	ompleted _	1040
Project N	lame	Limited	d Soil Invest	igation	Drill Rig			Geoprob	e 540 UD	
Geologist	t Jay F	erris	Driller	D. Riley	Auger Type		Macr	oCore 48-i	inch Direct	Push
Borehole	Completio	n/Abandon:	ment	Borehole allowe	i to collapse to s	urface	grade			
Depth (feet) Groundwater	Sample Type CS Symbol			,·		Recovery	PID Reading	Sample	Analytical	Sampling
a å	Symbol		Mate	erial Description		Rec	(ppm)	Collected	Method	Time
	SM	Dark browi	n, organic, sil	ty SAND (dry) (no	odor)				_	
1.0	4	<del> </del>		:		<b>////</b> _	<u></u>			
2.0										
,	СН	Dark brow (dry) (no o		orown, mottled, de	nse, plastic CLAY					
3.0										
			:							
4.0										
							·	· .	ļ	<u> </u>
5.0	- CH/SO	Brown to	reddish brow	n, mottled, sandy,	plastic CLAY (dr		<del> </del>		·	<u> </u>
		(10 000)						<u> </u>	· · · · ·	
6.0										
			·-·		·					
7.0		Tible -	to Parasi							
	SP	to heavy creosote								
8.0										
1	Sample T		oring 8.0 feet l	pelow surface grade Groundwate	<u> </u>	-		Co	mments	
E	Continuous Core First Detected Groundwater						*		: laboratory	analysis
1 =	Split Spoon Equilibrated Water						ppm:	parts per m		
_ i _	Hand Auger						NR:	no recover		

Project No.	C19	3.001 Fa	cility ID	ConSteelCo	Date Drilled	_	9/19/06	. To	tal Depth	8.0 ft.
ite1	625 Flow	ood Drive,	Flowood, M	Aississippi	Start Time	_	1045	Time C	ompleted _	1055
Project Nam	ie	Limited	Soil Invest	tigation	Drill Rig			Geoprob	e 540 UD	
Geologist _	Jay F	erris	Driller	D. Riley	Auger Type		Macr	oCore 48-	inch Direct	Push
Borebole Co	mpletion	/Abandonn	nent	Borehole allowed	l to collapse to s	urfa	ce grade			
Depth (feet) Groundwater Sample Type	USCS					Recovery	PID Reading	Sample	Analytical	Sampling
San Gre	Symbol		Mate	erial Description		lg Re	(ppm)	Collected	Method	Time
1.0						N R				,
1.0								<del> </del>		
2.0	SM	Dark brown	to black, si	ty fine SAND with	wood (railroad					<u> </u>
	SIVI	tie?) debris	(moist to we	t) (heavy creosote (	odor)					
3.0							·			
					· ·				\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	
4.0									-	
								<u> </u>		-
5.0	СН	Light gray		rm, plastic CLAY (	moist) (heavy				<u> </u>	·
								_		<del> </del>
6.0		<u> </u>	·							<u> </u>
7.0	SP	Dark gray odor)	to dark bro	wn, mottled, fine S	AND (moist) (hea	vy				
							<b></b>	<u>.</u>	_	10.50
8.0		Ì	·	· · · · · · · · · · · · · · · · · · ·				*B-7	82700	10:50
	Sample T		ring 8.0 feet	below surface grade  Groundwate					mments	
·			Dimet		<del></del>		*			malveie
	Continuous Core First Detected Groundward Split Spoon Equilibrated Water Le					*: ppm:	parts per m	r laboratory a nillion	marysis	
	, — · · · · · · · · · · · · · · · · · ·						NR:	no recover		

Project	t No.	C19	3.001	Facility:ID _	ConSteelCo	Date Drilled	_	9/19/06	To	tal Depth _	8.0 ft.
Site _	10	25 Flow	ood Drive	, Flowood, M	lississippi	Start Time		1055	Time C	ompleted _	1105
Projec	t Nam	e	Limite	d Soil Invest	igation	Drill Rig	_		Geoprob	e 540 UD	
Geolog	gist _	Jay F	erris	Driller	D. Riley	Auger Type		Macr	oCore 48-	inch Direc	Push
Boreh	ole Co	mpletion	/Abandon	ment	Borehole allowe	d to collapse to s	urfa	ce grade			
Depth (feet)	91 -1	USCS Symbol		Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
							N R				
1.0											· · · · · · · · · · · · · · · · · · ·
2.0		CL	Dark brow creosote o		plastic CLAY (m	oist) (slight					
3.0										<u> </u>	
				· · · · · ·	·					ļ.·	
4.0		CL/SP	Black, cre		soft, sandy CLAY	(moist) (heavy					
				:							
5.0											
		СН	Greenish odor)	gray, dense, p	lastic CLAY (moi	st) (heavy creosost	e ///				
6.0											
					·· · · · · · · · · · · · · · · · · · ·	, 				•	
7.0											
		SP		green, creosot reosote odor)	e-soaked/stained,	fine SAND (moist)					
8.0	)				-						
<b> </b>				oring 8.0 feet b	elow surface grade		T				_ <del></del>
	S	ample Ty	v <u>pe</u> uous Core	Rieet	Groundwat  Detected Groundw	<del></del>	<u>Comments</u> *: selected for laboratory analysis				analysis
Split Spoon Equilibrated Water Le							ppm:	parts per m	illion	anaiysis	
		Hand A	vuger					NR:	no recover	y 	·

Proj	ect ]	No.	<u>C193</u>	3.001 I	Facility ID	ConSteelCo	Date Drilled	_9	/19/06	Tot	tal Depth_	8.0 ft.
Site		1	625 Flow	ood Drive	, Flowood, M	ississippi	Start Time		1205	Time C	ompleted _	1215
Proj	ect	Nam	e	Limite	ed Soil Investi	gation	Drill Rig			Geoprob	e 540 UD	
Geo	logi	st _	Jay Fe	rris	Driller_	D. Riley	Auger Type		Macr	oCore 48-	inch Direc	Push
Bor			mpletion	/Abandon	ment <u>]</u>	Borehole allowed	to collapse to s	urfac	e grade			
Depth (feet)	Groundwater	Sample Type	USCS Symbol		Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
1.0	-		СН		vn, slightly silty noderate creosc	, firm, plastic CLA	aY (dry to moist)	N R		,	· .	
2.0	<u>,   .</u>			(sugnt to n	noderate creos(	ote odor)						
3.	0				· · · · · · · · · · · · · · · · · · ·		<u></u>					
4.	0		СМ		wn to black, sil eosote odor)	ty, soft, plastic CL	AY (moist to wet)					·
5	.0		The state of the s									
	5.0		СН	Greenish creosote		stic CLAY (dry to	moist) (heavy					
	7.0											
			CL/CN	1 Reddish	brown, friable	, slightly silty CLA	Y (heavy creosot					
	8.0		СН	<del></del>		astic CLAY (dry to	moist) (heavy					
			·		oring 8.0 feet b	elow surface grade		<u> </u>				
			Sample Ty Continu	r <u>pe</u> Jous Core	First	Groundwate Detected Groundwa			*:		mments laboratory	analysis
	Split Spoon Equilibrated Water Level ∇  Hand Auger						vel ∇		ppm: NR:	parts per m		

oject	No.	C19	3.001 F	facility ID	ConSteelCo	Date Drilled		9/19/06	To	tal Depth	8.0 ft.			
te	1	625 Flow	ood Drive,	, Flowood, N	<u> Iississippi</u>	Start Time		1215	Time C	ompleted _	1225			
roject	Nam	ie	Limite	d Soil Invest	igation	Drill Rig			Geoprob	e 540 UD	· · · · · · · · · · · · · · · · · · ·			
eolog	ist _	Jay F	erris	Driller	D. Riley	Auger Type		Macı	oCore 48-	inch Direc	t Push			
		mpletion	/Abandon	ment	Borehole allowed	l to collapse to s	urfac	e grade			·-····································			
Depth (feet)	Sample Type	USCS Symbol		Mate	erial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time			
						<del></del>								
.0				:.										
		CL	Dark gray		<del></del>									
2.0		CL		CLAY (dry)										
_						·								
3.0	_								Comments Selected for laboratory aparts per million					
_								·	-	<u> </u>	<u> </u>			
4.0					,				-		-			
_		СН	Gray, den	se, plastic CL	AY (dry to moist) (	ry to moist) (no odor)		· 		<del> </del>	<u> </u>			
5.0		MON HOUSE						·			-			
					•			· 		-	<u> </u>			
6.0				<del></del>	<del> </del>				-	+				
7.0		CL			reddish brown, mo									
			plastic C	LAY (dry) (m	oderate to heavy c	eosote odor)								
8.0			F 1 2=			<del></del>								
-		Sample T		oring 8.0 feet t	pelow surface grade Groundwate		+	<del> </del>	<u></u>	mmente				
			uous Core	Piret	Detected Groundwa	-		* •			analycic			
		Split S			quilibrated Water Le			ppm:						
		Hand A	Auger					NR:	no recover	у				

Projec	t No.	_C19	3.001	Facility:ID	ConSteelCo	Date Drilled	9	9/19/06	To	tal Depth_	8.0 ft.
Site _	1	625 Flow	ood Driv	e, Flowood, M	ississippi	Start Time		1230	Time C	ompleted _	1240
Projec	et Nam	ie	Limit	ed Soil Investi	gation	Drill Rig			Geoprob	e 540 UD	
Geolo	gist _	Jay F	erris	Driller_	D. Riley	Auger Type		Macr	oCore 48-	inch Direc	Push
		mpletion	ı/Abandoı	nment _	Borehole allowed	l to collapse to s	urfac	e grade			
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
1.0		CL/CM		ghtly silty to sil y) (no odor)	ty, slightly sandy,	low-plasticity					
SC Dark brown to black, silty, slightly clayey SAND with wood debris (dry to moist) (no odor)											
3.0 4.0 5.0		CL		y to dark reddi CLAY (dry) (r	sh brown, mottled o odor)	, friable, low-					
7.0		SP	Light gr	ray fine SAND (	moist) (no odor)						
		Com1 - 07		Boring 8.0 feet b	elow surface grade						
	Sample Type  Continuous Core  Split Spoon  Hand Auger  Groundwate  First Detected Groundwa  Equilibrated Water Le							* : ppm: NR:			analysis

Project No.	C19	3.001 F	acility ID	ConSteelCo	Date Drilled	9/19/06	То	tal Depth _	8.0 ft.
Site	1625 Flov	vood Drive,	Flowood, M	lississippi	Start Time	1245	Time C	Completed _	1255
Project Nai	ne	Limited	l Soil Investi	gation	Drill Rig		Geoprob	e 540 UD	
Geologist	Jay F	erris	Driller	D. Riley	Auger Type	Ma	croCore 48-	inch Direc	Push
Borehole C	ompletion	n/Abandonr	ment	Borehole allowed	d to collapse to s	urface grad	e		
Depth (feet) Groundwater Sample Type	uscs				·	PID Readin (ppm)	_	Analytical	
A G S	Symbol		Mate	rial Description		జ (ppm)	Collected	Method	Time
1.0	CL/CG		ı to black, silt l fill in upper	y, friable, low-plas 0.5 feet	sticity Clay with				
2.0									
3.0	CL		eddish brown /) (no odor)	mottled, friable, l	ow-plasticity				
4.0									
					·				
5.0									
6.0	СН	Dark gray (moist) (no		rown, mottled, den	se, plastic CLAY				-
7.0									
8.0			<u>.</u>						
	Sample T		oring 8.0 feet b	elow surface grade  Groundwate				mments	
		uous Core	First	Groundwate Detected Groundwa	<del></del>	*:		r laboratory	analysis
111	Split Spoon Equilibrated Water Lev					ppm NR			,

Proje	ct N	lo.	<u>C19</u>	3.001 F	acility ID	ConSteelCo	Date Drilled		9/19/06	To	tal Depth_	8.0 ft.
Site _		16	25 Flow	ood Drive,	Flowood, N	Aississippi	Start Time		1315	Time C	ompleted _	1325
Proje	ct N	Vamo	e	Limited	l Soil Invest	tigation	Drill Rig	_		Geoprob	e 540 UD	
Geolo	ogis	t _	Jay F	erris	Driller	D. Riley	Auger Type		Macr	Core 48-	inch Direct	Push
Borel	hole	e Co	mpletior	/Abandoni	ment	Borehole allowe	d to collapse to s	surfa	ce grade	<u>.</u>		
Depth (feet)	Groundwater	Sample Type	uscs					Recovery	PID Reading		Analytical	
ď	ق	Sa	Symbol		Mate	erial Description			(ppm)	Collected	Method	Time
1.0			SM/SC			ghtly silty to silty, s bris (moist) (no oc			·			
												· · · · · ·
2.0						, , , , , , , , , , , , , , , , , , ,			· · · · · · · · · · · · · · · · · · ·			
3.0	-		CL .			brown, mottled, fi	riable, low-					
4.0				plasticity (	CLAY (dry) (	no odor)						
-	-											
5.0	+											
6.0	D		SM	Black to recreosote o		n, silty, fine SAND	(moist) (slight					
	+				·							
7.			SP	Light bro	wn fine SAN	D (moist) (no odor	)					
8.	.0			End of Do	ring & O fact	below surface grade	· .					
		s	ample T		I GO TECT	Groundwat		-		Co	mments	
				uous Core		t Detected Groundw quilibrated Water L	ater <b>V</b>				r laboratory	analysis
	[	Ī	Hand A							no recover		

Projec	t No.	C19	3.001	Facility ID	ConSteelCo	Date Drilled	_9/	19/06	Tot	tal Depth_	8.0 ft.
Site _		1625 Flov	ood Driv	e, Flowood, N	Aississippi	Start Time		1330	Time C	ompleted _	1340
Proje	ct Nar	ne	Limit	ed Soil Invest	tigation	Drill Rig			Geoprob	e 540 UD	
Geolo	gist	Jay F	erris	Driller	D. Riley	Auger Type		Macr	oCore 48-	nch Direc	Push
Boreb	ole C	ompletio	ı/Abandoı	ıment	Borehole allowed	d to collapse to s	urface	grade			
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mate	erial Description			PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
1.0											·
2.0		CH/SC	(moist) (n	m, plastic, slig o odor)	thtly silty, slightly s	sandy CLAY					
4.0				· ·							
5.0		СН		y to reddish b ry to moist) (n	rown, mottled, den 10 odor)	se, highly plastic,					
7.0		SP	Light br	own, fine SAN	TD (moist) (no odor	)					
8.0			End of F	Boring 8.0 feet l	pelow surface grade					·	
Sample Type  Continuous Core  Split Spoon  Hand Auger  Groundwar  First Detected Groundwar  Equilibrated Water L						ater ▼		ppm:	-		analysis

roject N	o. <u>C</u> 1	93.001	Facility ID	ConSteelCo	Date Drilled	9/19	0/06	Tot	al Depth_	8.0 ft.
ite	1625 Flo	wood Driv	e, Flowood, N	<u> Iississippi</u>	Start Time	13	40	Time C	ompleted _	1350
roject N	ame	Limit	ted Soil Invest	igation	Drill Rig			Geoprob	e 540 UD	
eologist	Jay J	Ferris	Driller	D. Riley	Auger Type		Macro	Core 48-i	nch Direct	Push
orehole	Completio	n/Abando	nment	Borehole allowed	to collapse to s	urface g	rade	<u> </u>		
eet)	Type									
Depth (feet) Groundwater	e USCS Symbol		Mate	erial Description		g Res		Sample Collected	Analytical Method	Sampling Time
	CL		y to brown, slig CLAY with gra	htly silty to silty, fri	iable, low-					
1.0		ļ	<del></del>	<del></del>		<b>/////////////////////////////////////</b>				<del></del>
				•					,	
2.0					·			_		
3.0										
4.0	СН	Gray to		mottled, dense, pla	stic CLAY (dry)					
5.0										
								<del> </del>		
6.0										
										<u> </u>
7.0				<del></del>				·	-	·
	- SI	Light b	orown, fine SAN	VD (moist) (no odor)	· ·					
8.0										
			Boring 8.0 feet	below surface grade						
_	Sample '			Groundwate	<del>_</del>				mments	
1 =	Continuous Core First Detected Groundw								laboratory	analysis
1 -	Split Hand	∕el ∇	1 '	ppm: parts per million  NR: no recovery			٠			
	tiano	. rugoi					INIX. H	o recovery	·	·

Projec	et No.	C19	93.001 F	acility ID	ConSteelCo	Date Drilled	9/19	/06	Tot	al Depth_	8.0 ft.
Site _	1	625 Flov	vood Drive	, Flowood, I	Aississippi	Start Time	14	30	Time C	ompleted_	1440
Projec	ct Nan	1e	Limite	d Soil Inves	tigation	Drill Rig			Geoprob	e 540 UD	
Geolo	gist _	Jay F	erris	Driller	D. Riley	Auger Type		Macro	Core 48-i	nch Direct	Push
Boreb	ole C	mpletio	n/Abandon	ment	Borehole allowed	l to collapse to s	urface g	rade			
eet)	Water				· · · · · · · · · · · · · · · · · · ·						
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mat	erial Description		g Rea	ID ding pm)	Sample Collected	Analytical Method	Sampling Time
		СН	Dark brow	n to black, fir	m, silty CLAY (mo	ist) (no odor)			<del></del>		
1.0			ļ	<del></del>	<del></del>		<b>//</b>		. ' ·		
		Í									L
2.0			·								
		·		•							
3.0		,									
4.0		СН	Light gray (dry) (no c		rown, mottled, dens	se, plastic CLAY		•			
5.0				•	•						
6.0				÷		· ·					
7.0	)										
		SP	Light bro	own, fine SAN	ND (moist) (no odor						
8.0	0							·			
				oring 8.0 feet	below surface grade						
1		Sample T	·		Groundwate	-			-	mments	•
	Continuous Core First Detected Groundw						*: selected for laboratory analysis				analysis
	Split Spoon Equilibrated Water L. Hand Auger					vel ∇	1 -	-	arts per m		
	الت										

Projec	t No.	C1	93.001 F	acility ID	ConSteelCo	Date Drilled	_9	9/19/06	To	tal Depth_	8.0 ft.
Site _		1625 Flov	wood Drive,	Flowood, M	lississippi	Start Time		1445	Time C	ompleted _	1455
Projec	et Na	me	Limited	l Soil Invest	igation	Drill Rig			Geoprob	e 540 UD	
Geolo	gist	Jay F	erris	Driller	D. Riley	Auger Type	_	Macro	Core 48-	inch Direct	Push
Boreb	ole (	Completio	n/Abandoni	ment :	Borehole allowe	d to collapse to s	urfac	e grade			
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
				<del></del>							
1.0											
	_									<u> </u>	
2.0		СН		ldish brown,	dense, mottled, pla	stic Clay (dry)		· 			
3.0		Cn	(no odor)								
4.0					٠.			· · · · · · · · · · · · · · · · · · ·	-		
-	900000							· 	-		
5.0			ļ	<del></del> _	· · · · · · · · · · · · · · · · · · ·	- <u></u>				<u> </u>	ļ
_											
6.0											·
-		SP		rown with bla eosote odor)	ack striations, fine	SAND (moist)				<u> </u>	-
7.0											
8.0											
-		Sample T		oring 8.0 feet b	elow surface grade Groundwate	er		<u> </u>	Co	mments	
		n	nuous Core	First	Detected Groundwa	<del></del>		*:			analysis
	Split Spoon Equilibrated Water							ppm: parts per million			·
	III	Hand	Auger					NR:	no recover	Comments coted for laboratory analysis is per million	

Projec	t No.	C19	3.001 F	acility ID	ConSteelCo	Date Drilled	_	9/19/06	To	tal Depth _	8.0 ft.		
Site _	1	625 Flow	ood Drive,	Flowood, M	Iississippi	Start Time		1500	Time C	ompleted _	1510		
Projec	t Nan	1e	Limited	Soil Invest	igation	Drill Rig	_		Geoprob	e 540 UD			
Geolo	gist _	Jay Fe	erris	Driller	D. Riley	Auger Type		MacroCore 48-inch Direct Push					
Boreh		mpletion	/Abandonn	nent	Borehole allowe	d to collapse to s	urfa	ice grade		· · · · · ·			
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mate	erial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time		
1.0	-				:								
		СН			mottled, dense, pla	stic CLAY			-				
2.0			(moist) (ligh	nt to moderat	e creosote odor)					,			
3.0									<del> </del>				
					. <u></u>					·	·		
4.0											·		
-			·		•				<del>-   </del>	· ·			
5.0									<u> </u>		1		
6.0		SP			, mottled, slightly h) (moderate to str		r)						
7.0													
-				·	÷	.*							
8.0	0	麗	End of Bo	ring 8.0 feet l	pelow surface grade		_						
		Sample Ty		3.0.1001	Groundwat		$\dashv$		Co	mments			
1		Continu	uous Core	First	Detected Groundwa			*:	selected for	r laboratory	analysis		
		Split Sp	poon	Ed	quilibrated Water Le	evel V		ppm:	parts per m	nillion			
		Hand A	Auger					NR:	no recover	y			

Project	t No.	C19:	3.001	Facility ID	ConSteelCo	Date Drilled		9/19/06	То	tal Depth	8.0 ft.
Site				e, Flowood, M		Start Time	_	1530		ompleted	1540
				ed Soil Invest	<del></del>	Drill Rig	_			e 540 UD	
Geolog		Jay Fe		Driller	D. Riley	Auger Type		Macr		inch Direc	t Push
Boreh	ole Co	mpletion	/Abandor	ament .	Borehole allowe		urfa	ce grade			
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
					Dood specon			Фрил			
1.0					· ·						
2.0											
		СН		eddish brown, osote odor)	mottled, dense, pla	stic CLAY (dry)					
3.0											
5.0					÷	•					
				•				·			
4.0											
5.0											
		2000				•					
6.0		C.D.	*								
		SP SP	Light gr	ay, tine SAND	(dry) (faint creoso	te odor)					
7.0											
<u>                                   </u>						:					<del> </del>
8.0	<b>V</b>	SP	Light b	rown, fine SAN	D (moist to wet) (	ao odor)			-	-	
		Sample Ty		Boring 8.0 feet b	elow surface grade Groundwat	or			<u></u>	mments	
-			pe ious Core	Firet	Detected Groundwar	<del></del>		* •		r laboratory	analveis
1	55.53	Split Sp		- Y	uilibrated Water Le				parts per n		u.141.7313
		Hand A			minimum water Le			ppm: NR:	no recover		

# SOIL BORING LOG B-20 (group)

Group of shallow borings attempting to penetrate buried debris area

Projec	t No.	C19	3.001	Facility ID	ConSteelCo	Date Drilled		9/19/06	To	tal Depth	8.0 ft.
Site _	1	625 Flow	ood Dr	ive, Flowood	, Mississippi	Start Time		1600	Time C	ompleted _	1630
Projec	t Nam	e <u></u>	Lim	ited Soil Inv	estigation	Drill Rig			Geoprob	e 540 UD	
Geolog	gist _	Jay F	erris	Drill	er D. Riley	Auger Type	_	Macr	oCore 48-	nch Direc	Push
Boreb	ole Co	mpletion	/Aband	lonment	Borehole allowed	l to collapse to su	ırfa	ce grade	•		
Depth (feet)	Groundwater Sample Type	USCS Symbol		M	aterial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
1.0		СН			n, mottled, dense, pla	stic CLAY (dry)					
2.0			(no odo	r)							
3.0		· .		· · · · · · · · · · · · · · · · · · ·		· · · · · · · · · · · · · · · · · · ·		·			
4.0		Debris	(plywo	od, railroad ti	THIN 3.0-4.5 FEET E es) (debris typically w g beavy creosote odor	et with creosote	N R				
5.0									*B-20	8270C	11:20
6.0		СН			h brown, mottled, der or) (edge of debris are						
7.0											
8.0		SP			ND (moist) (no odor) (	edge of debris area	1)				
}	<del></del>	1 1 00		t Boring 8.0 fe	et below surface grade						
1		Sample Ty		_	Groundwate	<del></del>	-		-	mments	
		Continu Split S	uous Core	e F	irst Detected Groundwa Equilibrated Water Le			•	selected for parts per m	· laboratory :	analysis
		.Hand A			Equilibrated Watt Le	YOJ Y		• •	no recover		

Project	No.	C19	3.001 F	acility ID	ConSteelCo	Date Drilled		9/20/06	To	tal Depth	8.0 ft.	
Site	_ 1	.625 Flow	ood Drive	, Flowood, M	Iississippi	Start Time		0810	Time C	ompleted	0820	
 Project	t Nan	1e	Limite	d Soil Investi	igation	Drill Rig			Geoprob	e 540 UD		
Geolog	gist	Jay F	erris	Driller	D. Riley	Auger Type	_	MacroCore 48-inch Direct Push				
Boreho	ole C	mpletion	/Abandon	ment	Borehole allowe	d to collapse to s	surfa	ce grade				
Depth (feet)	Sample Type	USCS Symbol	``	Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time	
		SM	Dark brow	n, slightly sand	dy SILT (topsoil)	(moist) (no odor)						
1.0					·							
2.0		CL			orown, mottled, fr	iable, low-						
3.0			plasticity (	CLAÝ (dry) (n	o odor)	•		<del></del>			'	
3.0												
4.0				<del>,</del>	<del></del>							
5.0						•						
6,0		СН		enish gray to 1 LAY (moist) (1		ottled, dense, high	ly					
7.0	698											
8.0	100000		Tradition 1	0.06								
		Sample Ty		ornig o.v reet o	elow surface grade Groundwa		-		Co	mments	·	
	Continuous Core First Detected Groundway					ater <b>V</b>		*:		r laboratory	analysis	
	Split Spoon Equilibrated Water I Hand Auger					evel $ abla$		ppm: NR:	parts per m			

Projec	t No.	C1	93.001 F	acility ID	ConSteelCo	Date Drilled	_9	/20/06	To	tal Depth_	8.0 ft.
Site _		1625 Flov	vood Drive,	Flowood, N	<b>lississippi</b>	Start Time		0830	Time C	ompleted _	0840
Projec	ct Na	me	Limited	Soil Invest	igation	Drill Rig		·	Geoprob	e 540 UD	
Geolo	gist	Jay F	erris	Driller	D. Riley	Auger Type		Macr	oCore 48-	inch Direc	Push
Boreb	ole C	ompletio	n/Abandonr	nent	Borehole allowed	to collapse to s	urfac	e grade			
Depth (feet)	Groundwater Sample Type	USCS Symbol		Mate	erial Description			PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
		SM	Brown, silty	SAND with	roots (topsoil) (dry	(no odor)					
1.0											
2.0		ML.	Light brown	n to reddish	brown, mottled, fri	able, slightly				·	
3.0			sandy SILI	C (dry) (no od	or)	·					
4.0				·							
5.0						:					
6.0		СН		/ to reddish b Dist) (no odor	orown, mottled, high	nly plastic CLAY					
7.0			(dif to int	osty (uo ouoi							
8.0			End of Ro	oring 8.0 feet	pelow surface grade						
			ype wous Core	First	Groundwate Detected Groundwa	ter ▼		*:	selected for	mments	analysis
		Split S		E	quilibrated Water Le	vel V		ppm: NR:	no recover		

Projec	t No.	C19	3.001 F2	acility ID	ConSteelCo	Date Drilled	_9	/20/06	To	tal Depth_	8.0 ft.
Site _	1	625 Flow	ood Drive,	Flowood, N	Iississippi	Start Time		0840	Time C	ompleted _	0850
Projec	t Nan	ıe	Limited	Soil Invest	igation	Drill Rig			Geoprob	e 540 UD	
Geolog	gist	Jay F	erris	Driller	D. Riley	Auger Type	· 	Macro	Core 48-	inch Direc	Push
Boreh	ole C	mpletion	n/Abandonn	nent	Borehole allowed	l to collapse to s	urfac	e grade			
Depth (feet)	Groundwater Sample Type	USCS Symbol	,	Mate	erial Description			PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
		ML	Dark brown	slightly sand	ly SILT with organ	iics (topsoil)					
1.0					·						
		CL	Light brown	ı to reddish l	brown, mottled, sti	ff, friable slightly				-	
2.0		CL.	plastic CLA					. <u> </u>			
3.0											
4.0		СН			rown, mottled, stiff	, highly plastic		· · · · · · · · · · · · · · · · · · ·			
5.0			CLAY (mo	ist) (no odor	)			·			<del> </del>
-								<del></del>		·	-
6.0				<del></del>		<del> </del>				-	
7.0		SC			orown, mottled, slig	thtly clayey, fine		·			
		7.	SAIND (mo	oist) (no odo	r <b>)</b>				-		-
8.0			End of Da	ring Q A fact L	olow mustosa ana 1	·					
	End of Boring 8.0 feet below surface grade  Sample Type  Groundw				Groundwate		$\dashv$		Co	mments	
					Detected Groundwa			*:		r laboratory	analysis
					quilibrated Water Le				parts per m	•	-
		Hand A	•			·			no recover		

Projec	t No	•	C19	3.001 F	acility ID	ConSteelCo	Date Drilled	9/2	0/06	To	tal Depth _	8.0 ft.
Site _		162	5 Flow	ood Drive,	Flowood, M	Iississippi	Start Time	0	855	Time C	ompleted_	0910
Projec	t Na	me		Limited	Soil Invest	igation	Drill Rig			Geoprob	e 540 UD	
Geolo	gist		Jay F	erris	Driller	D. Riley	Auger Type		Macr	Core 48-	inch Direc	Push
Boreh	ole (	Com	pletion	ı/Abandonn	nent	Borehole allowed	l to collapse to s	urface	grade		<del></del>	
Depth (feet)	Groundwater	U	JSCS /mbol		Mate	rial Description		g Re	PID eading ppm)	Sample Collected	Analytical Method	Sampling Time
			ML	Dark brown	, slightly san	dy Silt with organi	cs (topsoil)					
1.0					<u>-</u>							
2.0	- The state of the											
			CL		n to reddish l LAY (dry) (n	brown, mottled, fri 10 odor)	able, low-					
3.0	1000000		=;								-	
	100									<u> </u>		
4.0	1				<del></del>	· · · · · · · · · · · · · · · · · · ·						<u> </u>
_									<del></del>	<b>-</b>	<u> </u>	
5.0							. •		· 			
_			СН	Light gree	nish gray to : AY (moist) (:	reddish brown, mo	ttled, dense, bighl	y 🖳 _				-
6.0					(						-	-
-	+									-		
7.0				_								-
8.0	0 🔻		SP	Light gra	y, fine SAND	(wet) (no odor)						
	1 13701				oring 8.0 feet b	elow surface grade						
1		Sai	mple T	ype		<u>Groundwat</u>	er			Co	mments	
	Continuous Core First Detected Ground				Detected Groundwa	ater <b>V</b>		*:	selected fo	r laboratory	analysis	
	Split Spoon Equilibrated Water L					evel $\nabla$		ppm: NR:	parts per n			
	Hand Auger							<u>. ]</u>	INK.	TIO TCCOVCI	<i>y</i>	

Proje	ct No.	_C19	03.001 F	acility ID	ConSteelCo	Date Drilled	9	9/20/06	Tot	al Depth_	8.0 ft.
Site _		1625 Flov	vood Drive,	Flowood, I	Mississippi	Start Time		0910	Time C	ompleted _	0920
Proje	ct Nar	ne	Limited	Soil Inves	tigation	Drill Rig			Geoprob	e 540 UD	
Geole	ogist	Jay F	erris	Driller	D. Riley	Auger Type		Macr	oCore 48-	inch Direct	Push
Bore	hole C	ompletio	n/Abandoni	nent	Borehole allowe	d to collapse to	surfac	e grade	· 		
eet)	vater										-
Depth (feet)	Groundwater Sample Type	USCS Symbol	<u>.</u>	Mat	erial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
		ML	Dark brown debris (tops		idy SILT with orga	nics and glass					
			debi is (tops	on,			₩				
1.0											
					•	•					
$\vdash$			\						<del> </del>		
2.0					-				· ·		
一			Brown to re	eddish brow	n, mottled, friable,	low-plasticity		<del></del>	<del> </del>	<del>                                     </del>	-
3.0		CL		to moist) (n		io ii piaoticity					
			ļ							T	
<u> </u>	-				•					ļ	<u> </u>
4.0					•						
-	1-8					·		·		<del> </del>	<del> </del>
5.0								·		1	
							-{//				
$\vdash$									<del>- </del> -		<del>-</del>
6.0									i		
-		SP	Brown to	light brown	fine SAND (moist t	o wet) (no odor)		]	_	<del></del>	<del></del>
7.	0		•		•				-		
_								<b></b>			· ·
8.	.0   ▼			•							
				ring 8.0 feet	below surface grade			<del> </del>			
1	احتجا	Sample T			Groundwat	<del></del>				mments	
			uous Core		t Detected Groundw			*:		laboratory	analysis
	188) (111)	Split S Hand	-	]	quilibrated Water Lo	evel V		ppm:	parts per m		
		riand.						NR:	no recovery	y 	

Proje	ct No	C1	93.001	Facility ID	ConSteelCo	Date Drilled	9	9/20/06	То	tal Depth	8.0 ft.
Site	•	1625 Flo	wood Driv	- e, Flowood, N	1ississippi	Start Time		0945	. Time C	ompleted	0955
Proje	ect Na			ed Soil Invest	· · · · · · · · · · · · · · · · · · ·	Drill Rig				e 540 UD	
Geol	ogist	Jay I	erris	Driller	D. Riley	Auger Type		Macr	oCore 48-	inch Direc	t Push
Bore	hole (	Completio	n/Abandoi	nment	Borehole allowe	d to collapse to s	urfac	e grade	_		
Depth (feet)	Groundwater	USCS Symbol		Mate	rial Description		Recovery	PID Reading (ppm)	Sample Collected	Analytical Method	Sampling Time
1.0											
2.0		СН		y to reddish br ry) (no odor)	own, mottled, dens	e, highly plastic					
3.0											
4.0						•					
5.0	)		Light gra	ay to reddish b	rown, mottled, find	e SAND (dry) (no					
6.0	0	SP				·					
7.	0		faint cre	eosote odor (7.0	)-8.0 feet below gra	ade)					
8.	.0		End of E	Boring 8.0 feet b	elow surface grade						
		Sample T			Groundwate	er			Co	mments	
		Contir	uous Core	First	Detected Groundwa	ater <b>V</b>		*:	selected for	r laboratory	analysis
		Split S Hand	-	Ec	uilibrated Water Le	vel V			parts per m		
- 1		-							•		

Project No. C193.001 Facility ID ConSteelCo Date Drilled 9/20/00	6 Total Depth 8.0 ft.
Site 1625 Flowood Drive, Flowood, Mississippi Start Time 0950	Time Completed1005
Project Name Limited Soil Investigation Drill Rig	Geoprobe 540 UD
Geologist Jay Ferris Driller D. Riley Auger Type M:	acroCore 48-inch Direct Push
Borehole Completion/Abandonment Borehole allowed to collapse to surface grad	de
(feet) Type Type	
PID Readin Symbol Material Description (ppm	ng Sample Analytical Sampling
CH Gray to reddish brown, dense, highly plastic CLAY (moist) (faint odor)	
3.0	
Gray to reddish brown, dense, slightly sandy, highly plastic	
CH/SP CH/SP CLAY (moist) (faint odor)	
5.0	
SP Light gray to reddish brown, mottled, fine SAND (moist)	
6.0	
CH/SP Light brown, slightly sandy, plastic CLAY (moist) (faint	
7.0 CH/SP odor)	
SP Light gray, fine SAND (moist) (faint odor)	
End of Boring 8.0 feet below surface grade	
Sample Type Groundwater	Comments
Continuous Core First Detected Groundwater ▼ *:	selected for laboratory analysis
Split Spoon Equilibrated Water Level ∇ ppr	n: parts per million
Hand Auger NF	R: no recovery

Appendix B Laboratory Analytical Report



Tax I.D. 62-0814289

Est. 1970

Mr. Jay Ferris Earth Consulting Group, Inc. P.O. Box 1246

Madison, MS 39130

#### Report Summary

Friday September 29, 2006

Report Number: L262154 Samples Received: 09/22/06 Client Project: C193.001

Description: ConSteelCo Creosote Assessment

The analytical results in this report are based upon information supplied by you, the client, and are for your exclusive use. If you have any questions regarding this data package, please do not hesitate to all.

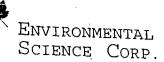
Reviewed By:

Craig Cothron, ESC Representative

#### Laboratory Certification Numbers

A2LA - 1461-01, AIHA - 09227, AL - 40660, CA - I-2327, CT - PH-0197, FL - E87487 GA - 923, IN - C-TN-01, KY - 90010, KYUST - 0016, NC - ENV375, DW21704, ND - R-140 NJ - TN002, SC - 84004, TN - 2006, VA - 00109, WV - 233 AZ - 0612, MN - 047-999-395, NY - 11742, WI - 998093910

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Tax I.D. 62-0814289

Est. 1970

REPORT OF ANALYSIS

Mr. Jay Ferris Earth Consulting Group, Inc. P.O. Box 1246 Madison, MS 39130

September 29, 2006

Date Received Description

September 22, 2006 ConSteelCo Creosote Assessment

ESC Sample # :

L262154-02

Sample ID

Site ID :

B-2 5.5 FT

Collected By : Collection Date :

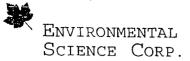
Jay Ferris 09/20/06 10:30

Project # :

C193.001

rameter	Result	Det. Limit	Units	Method	Date	n. i
se/Neutral Extractables					Date	Dil.
Acenaphthene						
Acenaphthylene	81.	66.	mg/kg	8270C	09/26/06	
Anthracene	BDL	66.	mg/kg	8270C		200
Benzidine	BDL	66.	mg/kg	8270C	09/26/06	200
Benzo(a) anthracene	BDL	66.	mg/kg	8270C	09/26/06	200
Benzo (b) fluoranthene	BDL	66.	mg/kg	8270C	09/26/06	200
Benzo(k) fluoranthene	BDL	66.	mg/kg	8270C	09/26/06	200
Benzo (g, h, i) perylene	EDL	66.	mg/kg	8270C	09/26/06	200
Benzo (a) pyrene	BDL	66.	mg/kg	8270C	09/26/06	200
Bis(2-chlorethoxy)methane	$\mathtt{BDL}$	66.	mg/kg	8270C	09/26/06	200
Bis(2-chloroethyl)ether	BDL	66.	mg/kg		09/26/06	200
Bis(2-chloroiser-	BDL	66.		8270C	09/26/06	200
Bis(2-chloroisopropyl)ether	BDL	66.	mg/kg	8270C	09/26/06	200
4-Bromophenyl-phenylether	BDL	66.	mg/kg	8270C	09/26/06	200
2-Chloronaphthalene	BDL	66.	mg/kg	8270C	09/26/06	200
4-Chlorophenyl-phenylether	BDL		mg/kg	8270C	09/26/06	200
Citysene	BDL	66.	mg/kg	8270C	09/26/06	200
Dibenz(a,h)anthracene	BDL	66.	mg/kg	8270C	09/26/06	200
3,3-Dichlorohenzidina		66.	mg/kg	8270C	09/26/06	
2,4-Dinitrotoluene	BDL	66.	mg/kg	8270C	09/26/06	200
2,6-Dinitrotoluene	BDL	66.	mg/kg	8270C		200
Fluoranthene	$\mathtt{BDL}$	66.	mg/kg	8270C	09/26/06	200
Fluorene	120	66.	mg/kg	8270C	09/26/06	200
Hexachlorobenzene	75.	66.	mg/kg	8270C	09/26/06	200
Hexachloro-1,3-butadiene	BDL	66.	mg/kg		09/26/06	200
Hexachlorogy-1-	BDL	66.		8270C	09/26/06	200
Hexachlorocyclopentadiene Hexachloroethane	BDL	66.	mg/kg	8270C	09/26/06	200
Indone	BDL	66.	mg/kg	8270C	09/26/06	200
Indeno(1,2,3-cd)pyrene	BDL	66.	mg/kg	8270C	09/26/06	200
Isophorone	BDL		mg/kg	8270C	09/26/06	200
Naphthalene	770	66.	mg/kg	8270C	09/26/06	200
Nitrobenzene	BDL	66.	mg/kg	8270C	09/26/06	200
n-Nitrosodimethylamine		66.	mg/kg	8270C	09/26/06	
N-Nitrosodiphenvlamine	BDL	10.	mg/kg	8270C	09/26/06	200
n-Nitrosodi-n-propylamine	BDL.	66.	mg/kg	8270C	09/26/06	200
richalthrene	BDL	66.	mg/kg	8270C		200
Benzylbutyl phthalate	200	66.	mg/kg	8270C	09/26/06	200
Bis(2-ethylhexyl)phthalate	$\mathtt{BDL}$	66.	mg/kg	8270C	09/26/06	500
Di-n-butyl phthalate	BDL	66.	mg/kg	8270C	09/26/06	200
Diethyl phthalate	BDL	66.	mg/kg		09/26/06	200
Dimethyl phthalate	BDL	66.		8270C	09/26/06	200
Dian control -late	BDL	66.	mg/kg	8270C	09/26/06	200
Di-n-octyl phthalate	BDL	66.	mg/kg	8270C	09/26/06	200
Pyrene	100		mg/kg	8270C	09/26/06	200
1,2,4-Trichlorobenzene	BDL	66.	mg/kg	8270C	09/26/06	200
id Extractables	801	66.	mg/kg	8270C	09/26/06	200

BDL - Below Detection Limit Det. Limit - Practical Quantitation Limit(PQL)



Tax I.D. 62-0814289

Est. 1970

REPORT OF ANALYSIS

Mr. Jay Ferris Earth Consulting Group, Inc. P.O. Box 1246 Madison, MS 39130

September 29, 2006

Date Received Description

September 22, 2006

ESC Sample # : L262154-02

ConSteelCo Creosote Assessment

Site ID :

Sample ID

Project # : C193.001

B-2 5.5 FT

Collected By : Collection Date :

Jay Ferris 09/20/06 10:30

Result	Det. Limit	Units	Method	Date	Dil.
BDL	66.	ma/ka	8270C	09/26/06	200
BDL					200
BDL					200
BDL					200
BDL					200
BDL					200
BDL					200
BDL.					200
					200
BDL					200
BDL					200
	•••	g/g	. 02100	03/20/00	200
0.00		t Rec	8270C	09/26/06	200
					200
			-		201
	•				20
					20
					20
	BDL BDL BDL BDL BDL BDL BDL BDL	BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66. BDL 66.	BDL 66. mg/kg BD	BDL 66. mg/kg 8270C BDL 66. mg/kg 8270C	BDL 66. mg/kg 8270C 09/26/06 BDL 66. mg/kg 8270C 09/26/06

BDL - Below Detection Limit

Det. Limit - Practical Quantitation Limit(PQL)

The reported analytical results relate only to the sample submitted.

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Tax I.D. 62-0814289

Est. 1970

REPORT OF ANALYSIS

Mr. Jay Ferris Earth Consulting Group, Inc. P.O. Box 1246 Madison, MS 39130

September 29, 2006

Date Received Description

September 22, 2006 ConSteelCo Creosote Assessment

ESC Sample # :

L262154-01

Sample ID

B-7 8 FT

Site ID :

Collected By : Collection Date :

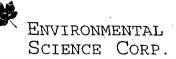
Jay Ferris 09/20/06 10:50

Project # :

C193.001

	Result	DetLimit	Units	Method	Date	
ase/Neutral Extractables					Date	Di.
Acenaphthene	200					
Acenaphthylene	390	66.	mg/kg	8270C	09/26/06	
Anthracene	BDL	66.	mg/kg	8270C		20
Benzidine	120	66.	mg/kg	8270C	09/26/06	20
Benzo(a) anthracene	BDL	66.	mg/kg	8270C	09/26/06	20
Benzo(b) fluoranthene	82.	66	mg/kg	8270C	09/26/06	. 20
Benzo(k) fluoranthene	BDL .	66.	mg/kg	8270C	09/26/06	20
Benzo(g,h,i)perylene	BDL	. 66.	mg/kg		09/26/06	20
Benzo(a) pyrene	$\mathtt{BDL}$	66.	mg/kg	8270C	09/26/06	20
Ric(2-oblomet	BDL	66.		8270C	09/26/06	20
Bis(2-chlorethoxy)methane	BDL	66.	mg/kg	8270C	09/26/06	20
Bis(2-chloroethyl)ether	BDL	66.	mg/kg	8270C	09/26/06	20
Bis(2-chloroisopropyl)ether	BDL		mg/kg	8270C	09/26/06	20
* "bromophenvi-phenvietha-	BDL	66.	mg/kg	8270C	09/26/06	20
<sup>2</sup> ~UllOronaphthalene	BDL	66.	mg/kg	8270C	09/26/06	20
4-Chlorophenyl-phenylether	BDL	66.	mg/kg	8270C	09/26/06	20
cut Aaeue		66.	mg/kg	8270C	09/26/06	20
Dibenz (a, h) anthracene	BDL	66.	mg/kg	8270C	09/26/06	
3,3-Dichlorobenzidine	BDL	66.	mg/kg	8270C		20
2,4-Dinitrotoluene	BDL	66.	mg/kg	8270C	09/26/06	20
2.6-Dinitrotoluene	BDL	66.	mg/kg	8270C	09/26/06	20
Fluoranthene	BDL	66.	mg/kg	8270C	09/26/06	20
Fluorene	490	66.	mg/kg		09/26/06	20
Hexachlorobenzene	370	66.	mg/kg	8270C	09/26/06	20
Hexachloro-1,3-butadiene	BDL	66.		8270C	09/26/06	20
Herachlono-1, 3-Ducadiene	BDL	66.	mg/kg	8270C	09/26/06	20
Hexachlorocyclopentadiene	BDL	66.	mg/kg	8270C	09/26/06	20
Hexachloroethane	BDL	66	mg/kg	8270C	09/26/06	20
Indeno(1,2,3-cd)pyrene	BDL	66.	mg/kg	8270C	09/26/06	20
Isophorone	BDL		mg/kg	8270C	09/26/06	20
Naphthalene	2300	66.	mg/kg	8270C	09/26/06	20
Nitrobenzene	BDL	660	mg/kg	8270C	09/29/06	20
n-Nitrosodimethylamine	BDL	66.	mg/kg	8270C	09/26/06	20
N-Nltrosodiphenvlamine		10.	mg/kg	8270C	09/26/06	
n-Nitrosodi-n-propylamine	BDL	66.	πg/)tg	8270C		20
Phenanthrene	BDL	66.	mg/kg	8270C	09/26/06	2(
Benzylbutyl phthalate	1000	330	mg/kg	8270C	09/26/06	2(
Bis(2-ethylhexyl)phthalate	BDL	66.	mg/kg	8270C	09/28/06	10
Di-n-butyl phthalate	BDL	66.	mg/kg		09/26/06	20
Diethyl phthalate	BDL	66.		8270C	09/26/06	20
Diethyl phthalate	BDL	66.	mg/kg	8270C	09/26/06	20
Dimethyl phthalate	BDL	66.	mg/kg	8270C	09/26/06	20
Di-n-octyl phthalate	BDL	66.	mg/kg	8270C	09/26/06	20
Pyrene	330		mg/kg	8270C	09/26/06	20
1,2,4-Trichlorobenzene	BDL	66.	mg/kg	8270C	09/26/06	20
id Extractables	מענו	66.	mg/kg	8270C	09/26/06	20

BDL - Below Detection Limit Det. Limit - Practical Quantitation Limit(PQL)



Tax I.D. 62-0814289

L262154-01

ßst. 1970

REPORT OF ANALYSIS

Mr. Jay Ferris Earth Consulting Group, Inc. P.O. Box 1246 Madison, MS 39130

September 29, 2006

Date Received

September 22, 2006

Description

ConSteelCo Creosote Assessment

Site ID :

ESC Sample # :

Sample ID

B-7 8 FT

Collected By : Collection Date :

Jay Ferris 09/20/06 10:50

Project # : C193.001

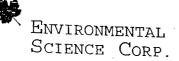
Parameter	Result	Det. Limit	Units_	Method	Date	Dil.
4-Chloro-3-methylphenol	BDL	66.	mg/kg	8270C	09/26/06	200
2-Chlorophenol	BDL	66.	mg/kg	8270C	09/26/06	200
2,4-Dichlorophenol	BDL	66.	mg/kg	8270C	09/26/06	200
2,4-Dimethylphenol	BDL	66.	mg/kg	8270C	09/26/06	200
4,6-Dinitro-2-methylphenol	BDL	66.	mg/kg	8270C	09/26/06	200
2,4-Dinitrophenol	BDL	66.	mg/kg	8270C	09/26/06	200
2-Nitrophenol	BDL	66.	mg/kg	8270C	09/26/06	200
4-Nitrophenol	· BDL	66.	mg/kg	8270C	09/26/06	200
Pentachlorophenol	BDL	66.	mg/kg	8270C	09/26/06	200
Phenol	BDL	66.	mq/kg	8270C	09/26/06	200
2,4,6-Trichlorophenol	BDL	66.	mg/kg	8270C	09/26/06	200
Surrogate Recovery				•	-	
Nitrobenzene-d5	0.00		* Rec.	8270C	09/26/06	200
2-Fluorobiphenyl	0.00		& Rec.	8270C	09/26/06	500
p-Terphenyl-dl4	0.00		* Rec.	8270C	09/26/06	200
Phenol-d5	0.00		* Rec.	8270C	09/26/06	200
2-Fluorophenol	0.00		₹ Rec.	8270C	09/26/06	200
2,4,6-Tribromophenol	0.00	•	* Rec.	8270C	09/26/06	200

BDL - Below Detection Limit Det. Limit - Practical Quantitation Limit(PQL) Note:

The reported analytical results relate only to the sample submitted.

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Tax I.D. 62-0814289

Est. 1970

REPORT OF ANALYSIS

Mr. Jay Ferris Earth Consulting Group, Inc. P.O. Box 1246 Madison, MS 39130

September 29, 2006

Date Received Description

September 22, 2006 ConSteelCo Creosote Assessment

ESC Sample # : L262154-03

Sample ID

Site ID :

B-20 5 FT

Collected By : Collection Date :

Jay Ferris 09/20/06 11:20

Project # : C193.001

arameter	Result	Det. Limit	Units	Method	Date	D 4 7
ase/Neutral Extractables					Date	Dil.
Acenaphthene	2400					
Acenaphthylene	2400	660	mg/kg	8270C	09/26/06	2000
Anthracene	BDL	660	mg/kg	8270C	09/26/06	2000
Benzidine	- 5500	660	mg/kg	8270C		2000
Benzo(a)anthracene	BDL	. 660	mg/kg	8270C	09/26/06	2000
Benzo (b) fluoranthene	930	660 -	mg/kg	8270C	09/26/06	2000
Benzo(k) fluoranthene	BDL	660	mg/kg	8270C	09/26/06	2000
Benzo (g, h, i) perylene	BDL	660	mg/kg	8270C	09/26/06	2000
Benzo (a) pyrene	BDL	660	mg/kg	8270C	09/26/06	2000
Bis(2-chlorethoxy)methane	$\mathtt{BDL}$	660	mg/kg		09/26/06	2000
Bis(2-chloroethyl)ether	$\mathtt{BDL}$	660	mg/kg	8270C	09/26/06	2000
Bis (2-chlorodernyl) ether	BDL	660		8270C	09/26/06	2000
Bis (2-chloroisopropyl) ether	BDL	660	mg/kg	8270C	09/26/06	2000
4-Bromophenyl-phenylether	BDL	660	mg/kg	8270C	09/26/06	2000
2-Chloronaphthalene	BDL	660	mg/kg	8270C	09/26/06	2006
4-Chlorophenyl-phenylether	BDL	660	mg/kg	8270C	09/26/06	2000
Cittysene	850		mg/kg	8270C	09/26/06	200
Dibenz (a, h) anthracene	BDL	660	mg/kg	8270C	09/26/06	2000
3,3-Dichlorobenzidine	BDL	660	mg/kg	8270C	09/26/06	200
2,4-Dinitrotoluene		660	mg/kg	8270C	09/26/06	200
2,6-Dinitrotoluene	BDL	660	mg/kg	8270C	09/26/06	
Fluoranthene	BDL	660	mg/kg	8270C	05/26/06	200
Fluorene	5400	660	mg/kg	8270C	09/26/06	200
Hexachlorobenzene	3300	660	mg/kg	8270C .	09/26/06	200
Hexachloro-1,3-butadiene	BDL	660	mg/kg	8270C	09/26/06	200
Hexachlorocyclopentadiene	BDL	660	mg/kg	8270C	09/26/06	200
Hexachloroethane	BDL	660	mg/kg		09/26/06	200
Indepo (1 2 2 - 1)	BDL	660		8270C	09/26/06	200
Indeno(1,2,3-cd)pyrene	BDL	660	mg/kg	8270C	09/26/06	200
Isophorone	BDL	660	mg/kg	8270C	09/26/06	200
Naphthalene	4600	660	mg/kg	8270C	09/26/06	200
Nitrobenzene	BDL		mg/kg	8270C	09/26/06	200
n-Nitrosodimethylamine	BDL	660	mg/kg	8270C	09/26/06	200
II-Nitrosodiphenvlamine	BDL	100	mg/kg	8270C	09/26/06	200
n-Nitrosodi-n-propylamine		660	mg/kg	8270C	09/26/06	200
rnenanthrene	BDL	660	mg/kg	8270C	09/26/06	
Benzylbutyl phthalate	7400	660	mg/kg	8270C	09/26/06	200
Bis(2-ethylhexyl) phthalata	BDL	660	mg/kg	8270C	09/26/06	200
Di-n-butyl phthalate	BDL	660	mg/kg	8270C	09/26/06	200
Diethyl phthalate	BDL	660	mg/kg		09/26/06	200
Dimethyl phthalate	$\mathtt{BDL}$	660	mg/kg	8270C	09/26/06	2000
Di-n-octyl phthalate	$\mathtt{BDL}$	660	mg/kg	8270C	09/26/06	200
Pyrene	BDL	660	mg/kg		09/26/06	200
1,2,4-Trichlorobenzene	3200	660	mg/kg	8270C	09/26/06	2000
cid Extractables	BDL	660		8270C	09/26/06	2000
DALLACEADIES		555	mg/kg	8270C	09/26/06	2000

BDL - Below Detection Limit
Det. Limit - Practical Quantitation Limit(PQL)

Page 5 of 8



Tax I.D. 62-0814289

Est. 1970

REPORT OF ANALYSIS

Mr. Jay Ferris Earth Consulting Group, Inc. P.O. Box 1246 Madison, MS 39130

September 29, 2006

Date Received September 22, 2006

ESC Sample # L262154-03

Description

ConSteelCo Creosote Assessment

Site ID :

Sample ID

B-20 5 FT

Collected By : Collection Date :

Jay Ferris 09/20/06 11:20

Project # : C193.001

Parameter	Result	Det. Limit	Units	Mahhad		
4-Chloro-3-methylphenol 2-Chlorophenol 2.4-Dichlorophenol 2.4-Dimethylphenol 4.6-Dimitro-2-methylphenol 2.4-Dinitrophenol 2.4-Dinitrophenol 2-Mitrophenol 4-Nitrophenol Pentachlorophenol Pentachlorophenol 2.4.6-Trichlorophenol Surrogate Recovery Nitrobenzene-d5 2-Fluorobiphenyl p-Terphenyl-d14 Phenol-d5 2-Fluorophenol 2.4.6-Tribromophenol	BDL BDL BDL BDL BDL BDL BDL BDL BDL BDL	660 660 660 660 660 660 660 660 660	mg/kg mg/kg mg/kg mg/kg mg/kg mg/kg mg/kg mg/kg mg/kg kg  Method  8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C 8270C	09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06 09/26/06	Dil.  2000 2000 2000 2000 2000 2000 2000 2	

BDL - Below Detection Limit
Det. Limit - Practical Quantitation Limit(PQL)
Note:

Note:
The reported analytical results relate only to the sample submitted.
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Reported: 09/29/06 16:05 Printed: 09/29/06 17:50

#### Attachment A List of Analytes with QC Qualifiers

Sample #	Analyte	Qualifíer
L262154-01	Acenaphthene	Ε
	Fluoranthene	E
•	Fluorene	E
	Naphthalene	E
	Phenanthrene	E
	Pyrene	E
	Nitrobenzene-d5	. J7
	2-Fluorobiphenyl	J7
	p-Terphenyl-d14	J7
	Pheno1-d5	J7
	2-Fluorophenol	J7
	2,4,6-Tribromophenol	່ 37
L262154-02	Naphthalene	E
	Phenanthrene	· E
	Nitrobenzene-d5	J7
	2-Fluorobiphenyl	J7
	p-Terphenyl-d14	J7 ·
	Phenol-d5	. 37
	2-Fluorophenol	J7
	2,4,6-Tribromophenol	J7
L262154-03	Acenaphthene	E
	Anthracene	· E
	Fluoranthene	E
•	Fluorene	E
	Naphthalene	E E E
	Phenanthrene	E
	Pyrene	E
	Nitrobenzene-d5	J7
	2-Fluorobiphenyl	J7
	p-Terphenyl-d14	J7
•	Phenol-d5	. <b>J7</b>
	2-Fluorophenol	J7
	2,4,6-Tribromophenol	37
•	• • • • • • • • • • • • • • • • • • • •	

#### Attachment B Explanation of QC Qualifier Codes

Qualifier	Meaning
E	GTL (EPA) - Greater than upper calibration limit: Actual value is known to be greater than the upper calibration range.
J7	Surrogate recovery limits cannot be evaluated; surrogates were diluted out
	Qualifier Report Information
as required by mos by ESC, we have im results. Each qua Data qualifiers ar the potential bias matrices incorpora	e and result qualifiers as set forth by the EPA Contract Laboratory Program and t certifying bodies including NELAC. In addition to the EPA qualifiers adopted plemented ESC qualifiers to provide more information pertaining to our analytical lifter is designated in the qualifier explanation as either EPA or ESC. e intended to provide the ESC client with more detailed information concerning of reported data. Because of the wide range of constituents and variety of ted by most EPA methods, it is common for some compounds to fall outside of these exceptions are evaluated and all reported data is valid and useable is 'R' (Rejected).
true va releva	Definitions lationship of the observed value of a known sample to the alue of a known sample. Represented by percent recovery and nt to samples such as: control samples, matrix spike recoveries, ate recoveries, etc.
Relat	greement between a set of samples or between duplicate samples. es to how close together the results are and is represented by ive Percent Differrence.
and c deter icall	ic compounds that are similar in chemical composition, extraction, hromotography to analytes of interest. The surrogates are used to mine the probable response of the group of analytes that are chemy related to the surrogate compound. Surrogates are added to the e and carried through all stages of preparation and analyses.  Control Limits (AO) (SS)
2-Fluoropheno Phenol-d5 2,4,6-Tribromophe	ol 31-119 Nitrobenzene-d5 43-118 Dibromfluoromethane 68-128 64-125 12-134 2-Fluorobiphenyl 45-128 Toluene-d8 76-115 69-118
notal	atively Identified Compound: Compounds detected in samples that are target compounds, internal standards, system monitoring compounds, urrogates.

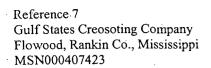
# Summary of Remarks For Samples Printed 09/29/06 at 17:50:25

TSR Signing Reports: 034 RX - Priority Rush

Sample: L262154-01 Account: ECGMS Received: 09/22/06 09:00 Due Date: 09/29/06 00:00 RPT Date: 09/29/06 16:05

Sample: L262154-02 Account: ECGMS Received: 09/22/06 09:00 Due Date: 09/29/06 00:00 RPT Date: 09/29/06 16:05

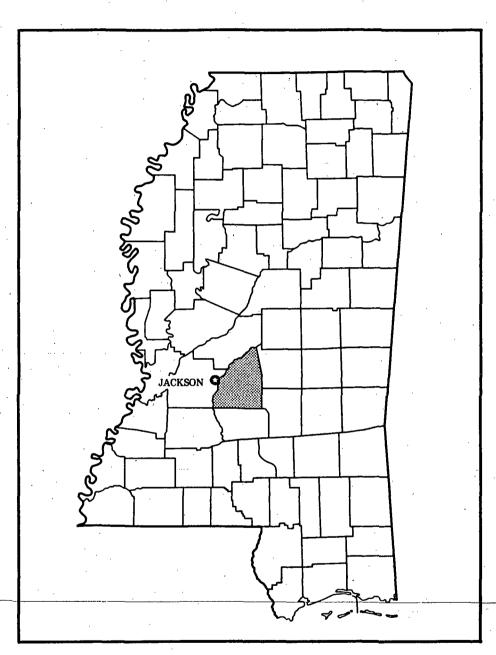
Sample: L262154-03 Account: ECGMS Received: 09/22/06 09:00 Due Date: 09/29/06 00:00 RPT Date: 09/29/06 16:05



with icultural xperiment Station

# Soil Survey of Rankin County, Ississippi





Location of Rankin County in Mississippi.

Water management (table 16)
Depth. USDA texture. Classification—Unified, AASHTO. Percentage passing sieve—4, 10, 40, 200. Liquid limit. Plasticity index.
Percentage passing sieve—4, 10, 40, 200. Liquid limit. Plasticity index.
Physical and chemical properties of the soils (table 18)
Depth. Clay. Moist bulk density. Permeability. Available water capacity. Soil reaction. Shrink-swell potential. Erosion factors. Organic matter.
Soil and water features (table 19)
Hydrologic group. Flooding. High water table. Risk of corrosion.
Physical and chemical analyses of selected soils (table 20) 177
Depth. Particle size distribution—Sand, Silt, Clay.
Extractable basis. Extractable acidity. Sums of catons. Base saturation. Reaction. Organic matter.
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5—Gillsburg silt loam, occasionally flooded6—Oaklimeter-Gillsburg association, frequently	19	42B—Providence-Urban land complex, 2 to 8 percent slopes
flooded	21	48C2—Ora fine sandy loam, 5 to 8 percent slopes,
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35D2-Tippah silt loam, 8 to 12 percent slopes,	35	64F—Smithdale-Providence association, hilly65D—Smithdale-Providence complex, 8 to 17
eroded	35	percent slopes66B—Providence-Tippah association, undulating
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	50	The state of the s

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# Soil Survey of Rankin County, Mississippi

By William A. Cole, Sr., Roger W. Smith, Mary Louise Spann, and Delmer C. Stamps, Soil Conservation Service

United States Department of Agriculture, Soil Conservation Service In cooperation with Mississippi Agricultural and Forestry Experiment Station

RANKIN COUNTY is in the southern part of Mississippi. It has a land area of 497,000 acres, or about 776.6 square miles. The total area, including bodies of water of more than 40 acres, is about 512,000 acres. Brandon, the county seat, is near the center of the county. The population of the county in 1980 was 68,183 according to the census.

The western boundary of the county is the Pearl River. The maximum dimension from north to south is about 37 miles and about 30 miles from east to west. The county is bounded on the north by Madison County, on the west by Hinds County, on the south by Simpson County, and on the east by Scott and Smith Counties.

Cotton, soybeans, forest products, poultry, beef and dairy production, and swine are the major sources of agricultural income in Rankin County. Many employees of nearby industrial plants are part-time farmers in the county.

The descriptions, names, and delineations of soils in this survey do not fully agree with those on soil maps for adjacent counties. Differences are the results of better understanding of soils, modification in series concepts, intensity of mapping, or the extent of soils within the survey area.

### General Nature of the Survey Area

This section provides information about the climate, history and development, transportation, physiography and geology, relief and drainage, and agriculture of Rankin County.

### Climate

Prepared by the National Climatic Data Center, Asheville, North Carolina.

Rankin County, Mississippi, has long, hot summers because moist tropical air from the Gulf of Mexico persistently covers the area. Winters are cool and fairly short with only a rare cold wave that moderates in 1 or days. Precipitation is fairly heavy throughout the year, and prolonged droughts are rare. Summer precipitation, mainly afternoon thunderstorms, is adequate for crops.

Table 1 gives data on temperature and precipitation for the survey area as recorded at Pelahatchie in the period 1951 to 1981. Table 2 shows probable dates of the first freeze in fall and the last freeze in spring. Table 3 provides data on length of the growing season.

In winter the average temperature is 50.3 degrees F, and the average daily minimum temperature is 37.7 degrees. The lowest temperature on record, which occurred at Pelahatchie on January 12, 1962, is -3 degrees. In summer the average temperature is 79 degrees, and the average daily maximum temperature is 91 degrees. The highest recorded temperature, which occurred on July 16, 1980, at Pelahatchie, is 104 degrees.

Growing degree days are shown in table 1. They are equivalent to "heat units." During the month, growing degree days accumulate by the amount that the averagatemperature each day exceeds a base temperature (50 degrees F). The normal monthly accumulation is used to schedule single or successive plantings of a crop between the last freeze in spring and the first freeze in fall.

The total annual precipitation is 55 inches. Of this, 26 inches, or 50 percent, usually falls in April through September. The growing season for most crops falls within this period. In 2 years out of 10, the rainfall in April through September is less than 21 inches. The heaviest 1-day rainfall during the period of record was 5.58 inches at Pelahatchie on December 29, 1954. Thunderstorms occur on about 64 days each year, and most occur in summer.

The average seasonal snowfall is less than 1 inch. The greatest snow depth at any one time during the period of record was 5 inches.

The average relative humidity in midafternoon is about 60 percent. Humidity is higher at night, and the average at dawn is about 90 percent. The sun shines 65 percent of the time possible in summer and 50 percent in winter. The prevailing wind is from the south. Average windspeed is highest, 9 miles per hour, in spring.

Severe local storms, including tornadoes, strike occasionally in or near the area. They are short and cause variable and spotty damage. Every few years in summer or autumn, a tropical depression or remnant of a hurricane that has moved inland causes extremely heavy rains for 1 to 3 days.

### **History and Development**

The earliest settlers in Rankin County were the Choctaw Indians. The Choctaw Indians farmed the land. In 1820, the Choctaws ceded to the United States the area that is now Rankin County in the Treaty of Doak's Stand, Second Choctaw Cecession (6).

On February 4, 1928, Rankin County was formed from the part of Hinds County that was east of the Pearl River. The county was named for Christopher Rankin, a congressman from Natchez and a member of the State Constitutional Convention of 1817. In 1829, commissioners were appointed to locate a site for the county seat. The site chosen was named Brandon in honor of Gerard C. Brandon, Mississippi's first native governor. In 1831, Brandon became the county seat. Early settlements in Rankin County included Richmond, Steens Creek, Fannin, Antioch, and Pisgah.

The Rankin County acreage that is west of Brandon and south to Florence is mainly in urban, commercial, and industrial use. This area adjoins the city of Jackson, Mississippi's state capitol. The remaining acreage is mainly in agricultural and woodland use.

The population of Rankin County in 1830 was 2,081, and in 1860, it had increased to 13,635. In 1979, the population was estimated at 65,000.

### **Transportation**

Access to Rankin County is provided by 10 state highways, two U.S. highways, Interstate Highways 55 and 20, and numerous county roads and streets. The Illinois Central Gulf Railroad has two lines that cross the

county and follow the U.S. highways in an east-west and northwest-south direction. Jackson Municipal Airport is in Rankin County and is serviced by three major airlines. Rankin County's proximity to Jackson, the state capitol, places it in the center of a busy transportation system.

### Physiography and Geology

Michael C. Seal, geologist, Mississippi Bureau of Geology, Jackson, Mississippi, prepared this section.

Mississippi is in the Gulf Coastal Plain physiographic province of North America. The state has been further subdivided into 12 physiographic units. In Rankin County, three of these units are represented. The northern two-thirds of the county is characterized by the Jackson Prairie Belt. Noted by gently rolling terrain, its southern limit roughly coincides with the geologic contact between the Yazoo Formation and the Forrest Hill Formation. South of the Jackson Prairie Belt is the Vicksburg Hills, characterized by gently rolling hills. The southern section of Rankin County is characterized by the Piney Woods physiographic unit and is underlain by the Catahoula Formation. On some of the higher elevations are outcrops of preloess terrace deposits and some Citronelle deposits (3).

Bedrock exposed in Rankin County is of Eccene, Oligocene, and Miccene series of the Tertiary System (4) and of the Pliestocene and Recent series of the Quaternary System.

The oldest unit exposed in the county is the Yazoo Formation of the Jackson Group. The Yazoo clay is a calcium montmorillonite that exhibits high shrink-swell potential with the removal or addition of water. This characteristic of the Yazoo clay causes foundation problems for all types of structures and roadbeds located on its outcrop. The major economic value of this material is as a lightweight aggregate, but it can also be mixed with other clays to make brick and ceramic materials.

The next oldest sediments exposed are of the Forrest Hill Formation. On the surface, the Forrest Hill sediments are thinly bedded, silty, micaceous, gray, fine to very fine grained sands. Clays are generally gray, buff, pink, and yellow. Thin lignite beds can also be observed in some outcrops. Petrified wood is often scattered over the surface of many Forrest Hill outcrops. A few domestic water wells are completed in the Forrest Hill Formation.

The Mint Springs marl is a gray-green, fine to coarse grained, fossiliferous to very fossiliferous, glauconitic sand. It is sometimes clayey in part and often limy. Fossils in this formation are mostly *Pectins* and oysters. Ferruginous sandstone, limonitic and manganiferous nodules and concretions are the end product of the weathered Mint Spring marl.

The Glendon limestone are characterized by alternating beds of limestone and mari. The thickness of the limestone beds ranges from 4 feet to less than 1/2

foot. They appear gray in the unweathered state. They are glauconitic, fossiliferous, and occasionally slightly sandy to sandy. Weathered Glendon limestone outcrops are noted by resistant limestone ledges, often intermittently apparent in dark-brown residuum. Some outcrops exhibit a white, waxy clay that is on the surface of the residuum. This clay is predominantly montmorillonite and halloysite with kaolinite as a trace—constituent. The Glendon limestone has produced lime that is suitable for agricultural and construction purposes. Many specimens of *Foraminifera* and *Pectins* can be collected at Glendon limestone outcrops.

Weathered Byram marl appears as brownish-red, slightly sandy clay. Ferruginous concretions are generally on the surface of weathered exposures. Fossils are abundant in the Byram marl, and several studies of these fossils have been made.

Weathered Bucatunna clay is chocolate brown, has conchoidal blocky fracture, is slightly micaceous, and is slightly silty. Some weathered Bucatunna clay has the resemblance of silty loam. Weathering can make Bucatunna clay difficult to identify.

In weathered outcrops, the Catahoula Formation is an indurated nonmarine series of clays, silts, and sands. In the unweathered state, it is generally not indurated. In some intervals of the Catahoula Formation in Rankin County, an extraordinary amount of salt is evident. Often, these have been used as salt licks by wild and domestic animals. The Catahoula Formation is a source of water for numerous domestic, agricultural, and municipal wells.

The Citronelle Formation is composed of chert and quartz gravel and fine to coarse grained sands and is Pliocene in age.

Preloess terrace deposits consist of fine- to coarsegrained sands that locally contain small amounts of pebble-size gravel. The sands are generally stained orange-red to buff and the gravels are generally finer than those in the Citronelle Formation.

Alluvial plains have developed along the two major rivers in Rankin County and along some of their tributaries.

### Relief and Drainage

The topography of Rankin County ranges from gently rolling to steep. In the north one-third of the county and in the area around the Jackson Dome, broad, rounded hills and wide, flat alluvial plains are common. In other parts, more sloping terrain is common. Some areas have high narrow ridges and deep narrow valleys. The highest elevation is about 612 feet along a ridge south of Shiloh Lookout Tower. The lowest point is in the southwest corner of the county along the Pearl River where the elevation is less than 220 feet.

Rankin County is drained by the Pearl River and its tributary, the Strong River. A ridge dividing the two watersheds crosses the southeastern one-fourth of the county in a southwest-northeast direction. Both rivers are fed by four main creeks and their tributaries. The major creeks in the Pearl River watershed are Fannegusha Creek in north Rankin County, Pelahatchie Creek in the north-central area, Richland Creek in west-central Rankin County, and Steen Creek in the southwestern area. The major creeks feeding the Strong River are the Dabbs, Campbell, Brushy, and Purvis Creeks.

### Agriculture

When Rankin County was inhabited mainly by the Choctaw Indians, corn was the major agricultural crop. Beans, pumpkins, and melons were the minor crops.

With the early European settlers came changing cropping systems, and before long, cotton was the major cash crop. About 7,500 bales of cotton were produced in 1851 and about 15,000 bales in 1899. Cotton production has fluctuated in the 20th century. About 4,500 bales were produced in 1924, 6,300 bales in 1969, 11,300 bales in 1974, and 9,200 bales in 1981. In recent years, poultry and poultry products have replaced cotton as the main cash crop. In 1974, cotton produced a total income of 2.5 million dollars while poultry and poultry products yielded more than 21.4 million dollars for Rankin County farmers.

Since the early 1900's, the number of farms in Rankin County has declined while the size of the farms has increased. There were 4,151 farms in 1910, 2,207 farms in 1925, and only 888 farms in 1974. The size of the average farm from 1910 to 1925 was 85 acres; and in 1974, it had increased to about 203 acres. In 1910, about 70 percent of the county was in farms, but by 1974, only 36 percent remained in farmland.

In 1965, about 3,000 dairy cattle were in Rankin County. By 1970 the number had declined to 1,700 and by 1974 it had declined to only 998. During this same period, the number of beef cattle changed little, and in 1974, it remained at about 23,000 head.

Woodland in Rankin County decreased from 359,900 acres in 1958 to 310,000 acres in 1977.

### **How This Survey Was Made**

This survey was made to provide information about the soils in the survey area. The information includes a description of the soils and their location and a discussion of the suitability, limitations, and management of the soils for specified uses. Soil scientists observed the steepness, length, and shape of slopes; the general pattern of drainage; the kinds of crops and native plants growing on the soils; and the kinds of bedrock. They dug many holes to study the soil profile, which is the sequence of natural layers, or horizons, in a soil. The profile extends from the surface down into the unconsolidated material from which the soil formed. The

unansolidated material is devoid of roots and other living aganisms and has not been changed by other biographical activity.

that is elated to the geology, the landforms, relief, climes and the natural vegetation of the area. Each kind is associated with a particular kind of landscape or with a segment of the landscape. By obsering the soils in the survey area and relating their position to specific segments of the landscape, a soil scients develops a concept, or model, of how the soils were timed. Thus, during mapping, this model enables the six scientist to predict with considerable accuracy the time of soil at a specific location on the landscape.

int: The another as their characteristics gradually charge. To construct an accurate soil map, however, soil scients must determine the boundaries between the soil. They can observe only a limited number of soil profes. Nevertheless, these observations, supplemented by an inderstanding of the soil-landscape relationship, are efficient to verify predictions of the kinds of soil in an efficient to determine the boundaries.

Spentists recorded the characteristics of the soil profes that they studied. They noted soil color, texture. size are shape of soil aggregates, kind and amount of roo: magments, distribution of plant roots, acidity, and other satures that enable them to identify soils. After describing the soils in the survey area and determining ther properties, the soil scientists assigned the soils to taxzzmic classes (units). Taxonomic classes are corrects. Each taxonomic class has a set of soil chazzeristics with precisely defined limits. The classes are see as a basis for comparison to classify soils system of taxonomic classification use: - the United States is based mainly on the kind and paracter of soil properties and the arrangement of homes within the profile. After the soil scientists classed and named the soils in the survey area, they concared the individual soils with similar soils in the sare axonomic class in other areas so that they could commutate and assemble additional data based on extended and research.

where a soil survey is in progress, samples of some of the sais in the area are generally collected for laboratory analyses and for engineering tests. Soil scientists interested the data from these analyses and tests as well as the field-observed characteristics and the soil processes in terms of expected behavior of the soils untal different uses. Interpretations for all of the soils were field tested through observation of the soils in different uses under different levels of management. Some interpretations are modified to fit local conditions, and the interpretations sometimes are developed to meetical needs. Data were assembled from other sources, such as research information, production recause, and field experience of specialists. For example,

data on crop yields under defined levels of management were assembled from farm records and from field or plot experiments on the same kinds of soil.

Predictions about soil behavior are based not only on soil properties but also on such variables as climate and biological activity. Soil conditions are predictable over long periods of time, but they are not predictable from year to year. For example, soil scientists can state with a fairly high degree of probability that a given soil will have a high water table within certain depths in most years, but they cannot assure that a high water table will always be at a specific level in the soil on a specific date.

After soil scientists located and identified the significant natural bodies of soil in the survey area, they drew the boundaries of these bodies on aerial photographs and identified each as a specific map unit. Aerial photographs show trees, buildings, fields, roads, and rivers, all of which help in locating boundaries accurately.

### Map Unit Composition

A map unit delineation on a soil map represents an area dominated by one major kind of soil or an area dominated by several kinds of soil. A map unit is identified and named according to the taxonomic classification of the dominant soil or soils. Within a taxonomic class there are precisely defined limits for the properties of the soils. On the landscape, however, the soils are natural objects. In common with other natural objects, they have a characteristic variability in their properties. Thus, the range of some observed properties may extend beyond the limits defined for a taxonomic class. Areas of soils of a single taxonomic class rarely, if ever, can be mapped without including areas of soils of other taxonomic classes. Consequently, every map unit is made up of the soil or soils for which it is named and some soils that belong to other taxonomic classes. In the detailed soil map units, these latter soils are called inclusions or included soils. In the general soil map units, they are called soils of minor extent.

Most inclusions have properties and behavioral patterns similar to those of the dominant soil or soils in the map unit, and thus they do not affect use and management. These are called noncontrasting (similar) inclusions. They may or may not be mentioned in the map unit descriptions. Other inclusions, however, have properties and behavior divergent enough to affect use or require different management. These are contrasting (dissimilar) inclusions. They generally occupy small areas and cannot be shown separately on the soil maps because of the scale used in mapping. The inclusions of contrasting soils are mentioned in the map unit descriptions. A few inclusions may not have been observed, and consequently are not mentioned in the descriptions, especially where the soil pattern was so

complex that it was impractical to make enough observations to identify all of the kinds of soils on the landscape.

The presence of inclusions in a map unit in no way diminishes the usefulness or accuracy of the soil data. The objective of soil mapping is not to delineate pure taxonomic classes of soils but rather to separate the

landscape into segments that have similar use and management requirements. The delineation of such landscape segments on the map provides sufficient information for the development of resource plans, but onsite investigation is needed to plan for intensive uses in small areas.

### **General Soil Map Units**

The general soil map at the back of this publication shows broad areas that have a distinctive pattern of soils, relief, and drainage. Each map unit on the general soil map is a unique natural landscape. Typically, a map unit consists of one or more major soils and some minor soils. It is named for the major soils. The soils making up one unit can occur in other units but in a different pattern.

The general soil map can be used to compare the suitability of large areas for general land uses. Areas of suitable soils can be identified on the map. Likewise, areas where the soils are not suitable can be identified.

Because of its small scale, the map is not suitable for planning the management of a farm or field or for selecting a site for a road or a building or other structure. The soils in any one map unit differ from place to place in slope, depth, drainage, and other characteristics that affect management.

Each map unit is rated for *cultivated crops, woodland, urban uses,* and *wildlife habitat*. Cultivated crops are those grown extensively in the survey area. Woodland refers to areas of native or introduced trees. Urban uses

include residential, commercial, and industrial developments. Wildlife habitat includes openland, woodland, and wetland wildlife habitat.

Dominantly nearly level soils that are well drained to poorly drained; on low stream terraces and flood plains

In this group are five general soil map units. The major soils are the well drained to poorly drained, silty Arkabutla, Cascilla, Gillsburg, Guyton, Leverett, Oaklimeter, Tippo, and Urbo soils; and the moderately well drained, loamy Kirkville and Quitman soils. The slopes range from 0 to 2 percent. These map units make up about 21.7 percent of the county.

### 1. Tippo-Leverett-Guyton

Nearly level, somewhat poorly drained, well drained, and poorly drained, silty soils; on low stream terraces and flood plains

This map unit consists of two broad areas in the westcentral part of Rankin County. These soils are on low

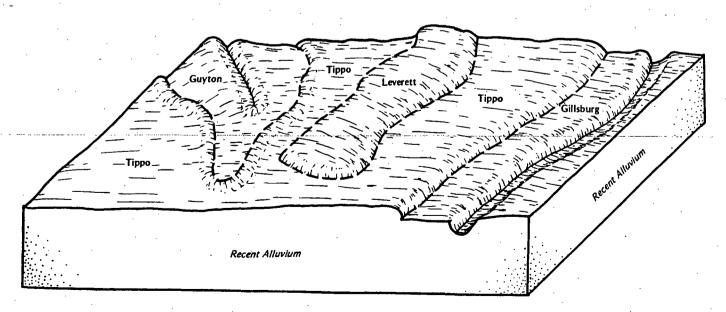


Figure 1.—The relationship of soils and landscape in the Tippo-Leverett-Guyton map unit.

stream terraces and flood plains along the Pearl River and its tributaries. The low stream terraces and flood plains have nearly linear surfaces. The topography is characterized by broad flats, low ridges, shallow swales, and winding stream channels (fig. 1). The soils in the low-lying areas are flooded after a heavy rain. Some depressions, swales, and sloughs are ponded during periods of unusual wetness. The slope ranges from 0 to 2 percent.

This map unit makes up about 3.2 percent of the county. It is about 50 percent Tippo soils, 15 percent Leverett soils, 14 percent Guyton soils, and 21 percent soils of minor extent.

Tippo soils are somewhat poorly drained and are on low stream terraces and flood plains. These soils formed in silty material. Leverett soils are well drained, are on low stream terraces, and are in slightly higher positions on the landscape than Tippo and Guyton soils. These soils formed in silty material. Guyton soils are poorly drained and are on broad, wet flats, stream terraces, and flood plains. These soils formed in silty alluvium.

The minor soils in this map unit are Gillsburg and Oaklimeter soils. These soils are silty, and they are on the flood plains. Gillsburg soils are somewhat poorly drained. Oaklimeter soils are moderately well drained.

The soils in this map unit are mostly in the urban areas of Flowood, Pearl, and Richland. In some areas, the soils are used for crops and pasture. Low, wet areas are in bottom land hardwoods.

Tippo and Leverett soils are well suited to row crops and small grains and to pasture grasses and legumes. Guyton soils are well suited to pasture grasses and legumes but are poorly suited to row crops because of wetness and flooding.

Tippo and Guyton soils are well suited to use as woodland, and Leverett soils are moderately suited to this use. Concerns in management are slight for use of Leverett soil as woodland. Wetness and flooding are severe limitations for use of equipment on Guyton soils and are moderate limitations on Tippo soils.

Guyton and Tippo soils have severe limitations for urban use because of flooding and wetness. In areas that are protected from flooding by levees, Tippo soils are moderately suited to urban use, and Guyton soils are poorly suited to this use because of wetness. Leverett soils have slight limitations for many urban uses.

Leverett and Tippo soils have good potential for the development of habitat for openland and woodland wildlife. Guyton soils have fair potential. For the development of habitat for wetland wildlife, Tippo soils have fair potential, Leverett soils have poor potential, and Guyton soils have good potential.

### 2. Cascilla-Arkabutla

Nearly level, well drained and somewhat poorly drained, silty soils; on flood plains

This map unit is in the western and northern parts of Rankin County. These soils mainly are on the flood plains of the Pearl River and its tributaries. The nearly linear surface of the flood plain is broken at irregular intervals by old river runs, natural levees, sloughs, chutes, and scarps (fig. 2). The slope ranges from 0 to 2 percent.

This map unit makes up about 3.7 percent of the county. It is about 40 percent Cascilla soils, 32 percent Arkabutla soils, and 28 percent soils of minor extent.

Cascilla soils are well drained. They are near the low scarps and on the slightly higher elevations on natural levees on flood plains along the Pearl River and the major tributaries. These soils formed in silty alluvium. Arkabutla soils are somewhat poorly drained. They are in broad, level areas, in slight depressions, and in the main flood basins of the flood plain. These soils formed in silty alluvium.

The minor soils in this map unit are the Gillsburg and Oaklimeter soils. These soils are silty and on the flood plains. Gillsburg soils are somewhat poorly drained. Oaklimeter soils are moderately well drained.

Most of the acreage in this map unit is in woodland. Because of wetness and flooding, Cascilla and Arkabutla soils are poorly suited to row crops and small grains. They are moderately suited to pasture grasses and legumes.

The soils in this map unit are well suited to use as woodland. Productivity is high for bottom land hardwoods. The use of equipment is limited because of wetness and flooding. Seedling mortality and plant competition are moderate limitations on these soils.

The soils in this map unit have severe limitations for urban use because of flooding.

Cascilla and Arkabutla soils have fair potential for development of habitat for openland wildlife and good potential for development of habitat for woodland wildlife. For development of habitat for wetland wildlife, Cascilla soils have very poor potential and Arkabutla soils have fair potential.

### 3. Urbo-Arkabutla

Nearly level, somewhat poorly drained, silty soils; on flood plains

This map unit is in the northern and north-central parts of Rankin County. These soils are along Pelahatchie and Fannegusha Creeks and their tributaries. Areas of these soils are subject to occasional or frequent flooding generally during winter or early in the spring. The slopes range from 0 to 2 percent.

This map unit makes up about 3.5 percent of the county. It is about 48 percent Urbo soils, 28 percent Arkabutla soils, and 24 percent soils of minor extent.

Urbo soils are on broad flats and in depressions of flood plains. These soils formed in clayey alluvium.

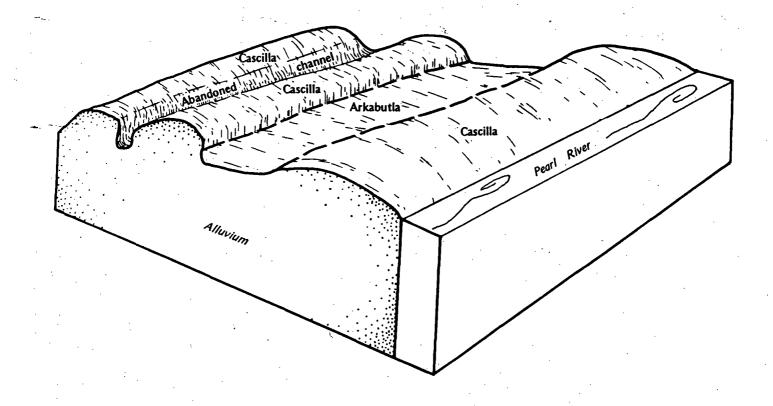


Figure 2.—The relationship of soils and landscape in the Cascilla-Arkabutia map unit.

Arkabutla soils are on broad flats of flood plains. These soils formed in silty alluvium.

The minor soils in this map unit are Gillsburg, Guyton, and Quitman soils. Gillsburg soils are somewhat poorly drained and are on the flood plains. Guyton soils are poorly drained and are on broad flats, stream terraces, and flood plains. Quitman soils are moderately well drained and are on stream terraces.

Most areas of this map unit are used as woodland.

Areas of this map unit that are frequently flooded are poorly suited to row crops and small grains and are only moderately suited to grasses and legumes. Areas that are occasionally flooded are well suited to most commonly grown crops and to grasses and legumes.

Urbo and Arkabutla soils are well suited to use as woodland. Productivity is high for bottom land hardwoods. The use of equipment is limited because of wetness and flooding. Seedling mortality and plant competition are moderate limitations on these soils.

The soils in this map unit have severe limitations for urban use because of flooding.

Urbo and Arkabutla soils have fair potential for the development of habitat for openland wildlife and wetland wildlife. For the development of habitat for woodland

wildlife, Urbo soils have fair potential and Arkabutla soils have good potential.

### 4. Oaklimeter-Gillsburg

Nearly level, moderately well drained and somewhat poorly drained, silty soils; on flood plains

This map unit is on flood plains in the southwestern part of Rankin County. The flood plains range from 300 feet wide to about 2 miles or more wide. Flooding is occasional or frequent. The slope ranges from 0 to 2 percent.

This map unit makes up about 5.3 percent of the county. It is about 58 percent Oaklimeter soils, 24 percent Gillsburg soils, and 18 percent soils of minor extent.

Oaklimeter soils are moderately well drained. They commonly are in slightly higher positions on the flood plains than Gillsburg soils. These soils formed in silty alluvium. The Gillsburg soils are somewhat poorly drained. They are on flood plains. These soils formed in silty alluvium.

The minor soils in this map unit are Arkabutla, Kirkville, Guyton, and Tippo soils. The Arkabutla soils are somewhat poorly drained and are on the flood plains.

Kirkville soils are moderately well drained and are on flood plains. Guyton soils are poorly drained and are on broad flats, stream terraces, and flood plains. The Tippo soils are somewhat poorly drained and are on broad flats, stream terraces, and flood plains.

Most of the acres of this map unit is used as woodland. Some areas are used for pasture or crops.

Areas of this map unit that are occasionally flooded are well suited to cultivated crops and small grains and to pasture grasses and legumes. Areas that are subject to frequent flooding are poorly suited to row crops and small grains.

Oaklimeter and Gillsburg soils are well suited to use as woodland, especially bottom land hardwoods. Flooding and seasonal wetness are the main concerns in woodland management and limit the use of equipment on these soils. Plant competition and seedling mortality are moderate limitations.

The soils in this map unit have severe limitations for urban use because of flooding.

Oaklimeter and Gillsburg soils have fair potential for the development of habitat for openland wildlife and good potential for the development of habitat for woodland wildlife. For development of habitat for wetland wildlife, Oaklimeter soils have poor potential and Gillsburg soils have fair potential.

#### 5. Quitman-Kirkville

Nearly level, moderately well drained, loamy soils; on low stream terraces and flood plains

This map unit is in the central and southeastern part of Rankin county. These soils mostly are on terraces and flood plains along the Strong River and its major tributaries (fig. 3). Kirkville soils are subject to flooding mainly during winter or early in the spring. Flooding is generally of brief duration. The slope ranges from 0 to 5 percent.

This map unit makes up about 6 percent of the county. It is about 54 percent Quitman soils, 34 percent Kirkville soils, and 12 percent soils of minor extent.

Quitman soils are on low terraces or second bottoms adjacent to the uplands. These soils formed in loamy material. Kirkville soils are on flood plains near stream channels. These soils formed in loamy alluvium.

The minor soils in this map unit are Guyton, Tippo, Oaklimeter, and Savannah soils. Guyton and Tippo soils are on broad flats, stream terraces, and flood plains. Guyton soils are poorly drained, and Tippo soils are somewhat poorly drained. Oaklimeter soils are moderately well drained and are on flood plains. Savannah soils are moderately well drained and are on stream terraces.

The soils in this map unit are used mainly for cultivated crops or as woodland. The other soils are in pasture. These soils are well suited to most commonly grown crops and small grains and to pasture grasses and legumes.

Quitman and Kirkville soils are well suited to use as woodland. Flooding and wetness are moderate limitations to use of equipment. Plant competition is a moderate limitation.

The Quitman soils in this map unit are moderately suited to urban use because of wetness. The Kirkville soils are poorly suited to urban use because of flooding.

Quitman and Kirkville soils have good potential for the development of habitat for openland and woodland wildlife and poor potential for the development of habitat for wetland wildlife.

### Dominantly nearly level to steep soils that are well drained to somewhat poorly drained; on uplands and stream terraces

In this group are five general soil map units. The major soils are the somewhat poorly drained to moderately well drained, silty Falkner, Kipling, Providence, and Tippah soils; and the moderately well drained to well drained, loamy Quitman, Savannah, and Smithdale soils. The slopes range from 0 to 40 percent. These map units make up about 78.3 percent of the county.

### 6. Kipling-Falkner-Savannah

Nearly level to sloping soils; some are somewhat poorly drained, silty soils that are underlain by a plastic, clayey subsoil; and some are moderately well drained, loamy soils that have a fragipan; on uplands and stream terraces

This map unit is on the prairie in the northern part of Rankin County. The landscape has low relief and is mainly nearly level to gently rolling. In some places, the low hills have a cap of loamy terrace sediments (fig. 4). The slope ranges from 0 to 8 percent.

This map unit makes up about 23.6 percent of the county. It is about 40 percent Kipling soils, 18 percent Falkner soils, 16 percent Savannah soils, and 26 percent soils of minor extent.

Kipling soils are silty and are somewhat poorly drained. They are on uplands. These soils formed in clayey material. Falkner soils are silty and are somewhat poorly drained. They are on uplands and stream terraces. These soils formed in a silty mantle and the underlying acid, clayey deposits. Savannah soils are loamy and moderately well drained and have a fragipan. They are in slightly higher positions on the uplands and stream terraces than Kipling and Falkner soils. These soils formed in loamy material.

The minor soils in this map unit are Pelahatchie, Providence, Quitman, and Urbo soils. Pelahatchie soils are moderately well drained and are on uplands. Providence and Quitman soils are moderately well drained and are on uplands and stream terraces. Urbo soils are somewhat poorly drained and are on the flood plains.

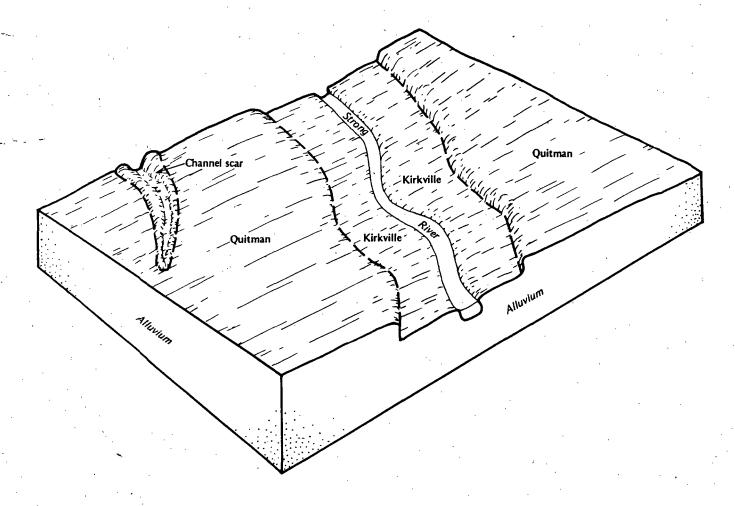


Figure 3.—The relationship of soils and landscape in the Quitman-Kirkville map unit.

Most of the acreage in this map unit is used for cultivated crops or as woodland. The other acreage is in pasture.

The nearly level or gently sloping areas of Falkner and Kipling soils are moderately suited to most commonly grown row crops and small grains and are well suited to grasses and legumes for hay and pasture. Savannah soils are well suited to row crops and small grains and to grasses and legumes for hay and pasture.

Kipling and Falkner soils are well suited to use as woodland. Seasonal wetness is a moderate limitation to use of equipment, and plant competition is a moderate limitation if pines are planted. Savannah soils are moderately suited to use as woodland, but windthrow and plant competition are moderate limitations.

Wetness and high shrink-swell potential of the subsoil severely restrict Kipling and Falkner soils for urban use. Mainly because of seasonal wetness, Savannah soils have moderate limitations for urban use. The soils in this map unit have good potential for the development of habitat for openland and woodland wildlife. For development of habitat for wetland wildlife, Falkner and Savannah soils have very poor potential. In the nearly level areas, Kipling soils have fair potential for habitat for wetland wildlife; in the gently sloping areas, they have poor potential; and in the sloping areas, they have very poor potential.

### 7. Smithdale-Providence

Gently sloping to steep soils; some are well drained, loamy soils; and some are moderately well drained, silty soils that have a fragipan; on uplands and steam terraces

This map unit is in the central and southern parts of Rankin County. The landscape is hilly and is marked by narrow ridgetops that are generally less than one-eighth of a mile wide, by hillsides that are dissected by many

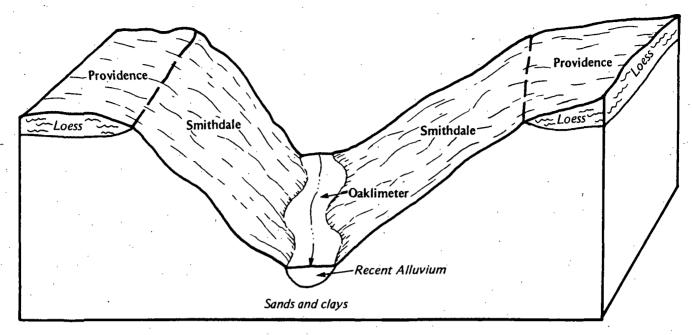


Figure 4.—The relationship of soils and landscape in the Kipling-Faikner-Savannah map unit.

short drainageways, and by narrow flood plains (fig. 5). The slope is dominantly 5 to 40 percent.

This map unit makes up about 19.3 percent of the county. It is about 43 percent Smithdale soils, 30 percent Providence soils, and 27 percent soils of minor extent.

Smithdale soils are loamy and are well drained. They are on the steeper hillsides on uplands. These soils formed in loamy material. Providence soils are silty and moderately well drained and have a fragipan. They are on uplands and stream terraces. These soils formed in a mantle of silty material and in the underlying loamy sediment.

The minor soils in this map unit are Kisatchie, Savannah, Tippah, Oaklimeter, Kirkville, and Gillsburg soils. Kisatchie soils are well drained and are on uplands. Savannah and Tippah soils are moderately well drained and are on uplands. Oaklimeter and Kirkville soils are moderately well drained and are on the flood plains. Gillsburg soils are somewhat poorly drained and are on the flood plains.

Most areas of this map unit are used as woodland. A small acreage is used for pasture and crops.

The Smithdale soils are poorly suited to row crops and small grains and to pasture grasses and legumes because of steep slopes. In the gently sloping areas, Providence soils are well suited to row crops, and in the sloping areas, they are moderately suited to this use. In the gently sloping and sloping areas, Providence soils are well suited to grasses and legumes for hay and

pasture, and in the sloping areas, they are moderately suited to this use.

Providence soils are moderately suited to use as woodland. Concerns in woodland management are few. Smithdale soils are moderately suited to woodland use. Steepness of slope is a moderate limitation to use of equipment on Smithdale soils if slopes are more than 15 percent.

Smithdale soils have severe limitations for urban use because of steepness of slopes. Providence soils have moderate limitations for urban use mainly because of seasonal wetness and steepness of slopes.

Smithdale and Providence soils have good potential for the development of habitat for openland and woodland wildlife, but on Smithdale soils if slopes are more than 15 percent, potential is fair. For the development of habitat for wetland wildlife, the potential of the soils in this map unit is very poor.

### 8. Providence-Tippah

Gently sloping to moderately steep, moderately well drained, silty soils; some have a fragipan; and some are underlain by plastic, clayey material; on uplands and stream terraces

This map unit is in the west-central and southwestern part of Rankin County. The landscape has moderate relief and is generally rolling but is moderately steep along the major drainageways. It is marked by broad ridgetops, by hillsides that are dissected by short

drainageways, and by narrow flood plains along the streams. The slope ranges from 0 to 15 percent.

This map unit makes up about 17.9 percent of the county. It is about 54 percent Providence soils, 31 percent Tippah soils, and 15 percent soils of minor extent.

Providence soils have a fragipan. They are on uplands and stream terraces. These soils formed in a mantle of silty material and the underlying loamy material. Tippah soils are on uplands. These soils formed in a mantle of silty material and the underlying clayey material.

The minor soils in this map unit are Kirkville, Oaklimeter, and Savannah soils. Kirkville and Oaklimeter soils are moderately well drained and are on the flood plains. Savannah soils are moderately well drained and are on uplands and stream terraces.

Most of the acreage in this map unit is used as woodland or pasture. Some areas are used for row crops.

In the gently sloping areas, Providence and Tippah soils are well suited to most commonly grown crops and small grains and to grasses and legumes for hay and pasture. In the sloping areas, these soils are moderately suited to most commonly grown crops and small grains and are well suited to grasses and legumes for hay and pasture.

The soils in this map unit are moderately suited to use as woodland. Plant competition is the main limitation on Tippah soils, and windthrow is a limitation on Providence soils.

Wetness and steepness of slopes are moderate limitations to use of Providence soils for urban use. On Tippah soils, wetness and shrink-swell potential of the subsoil are severe limitations for urban use.

The soils in this map unit have good potential for the development of habitat for openland and woodland wildlife and poor or very poor potential for development of habitat for wetland wildlife.

### 9. Smithdale-Savannah

Gently sloping to steep, loamy soils; some are well drained; and some are moderately well drained and have a fragipan; on uplands and stream terraces

This unit is in the eastern and southeastern parts of Rankin County. The landscape is hilly and is marked by narrow ridgetops that are generally less than one-eighth of a mile wide, by steep hillsides that are dissected by many short drainageways, and by narrow flood plains along the streams. The slope ranges from 2 to 40 percent.

This map unit makes up about 11.4 percent of the county. It is about 44 percent Smithdale soils, 38 percent Savannah soils, and 18 percent soils of minor extent.

Smithdale soils are well drained. They are on the steeper upland hillsides than Savannah soils. These soils

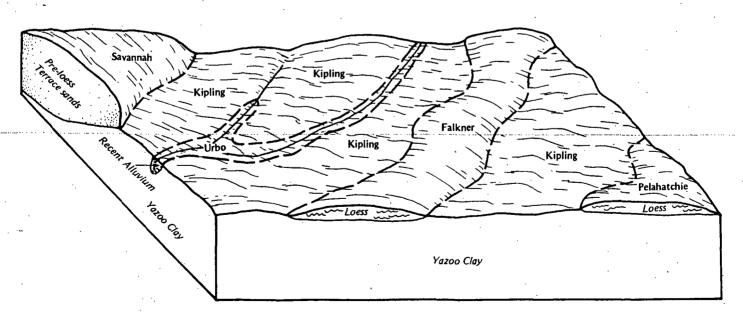


Figure 5.—The relationship of soils and landscape in the Smithdale-Providence map unit.

formed in loamy marine sediment. Savannah soils are moderately well drained and have a fragipan. They mainly are on ridgetops on uplands and stream terraces. These soils formed in loamy material.

The minor soils in this map unit are Kirkville, Maben, Ora, and Tippah soils. Kirkville soils are moderately well drained and are on the flood plains. Maben soils are well drained and are on uplands. Ora and Tippah soils are moderately well drained and are on uplands.

Most of the acreage in this map unit is used as woodland or pasture. A small acreage is in row crops.

Smithdale soils are poorly suited to row crops and small grains but are moderately suited to grasses and legumes. The main limitation to use of these soils for row crops and small grains is the steep slopes. Erosion is a hazard. In the gently sloping areas, Savannah soils are well suited to row crops and small grains; in the sloping areas, they are moderately suited to this use. Savannah soils are well suited to grasses and legumes for hay and pasture.

The soils in this map unit are moderately suited to use as woodland. Smithdale soils have no significant limitations to woodland use and management. Plant competition and windthrow are moderate limitations to use of Savannah soils as woodland if pine trees are planted.

Smithdale soils have severe limitations for urban and recreational uses because of steepness of slope. Savannah soils are moderately limited for these uses because of wetness.

Smithdale and Savannah soils have good potential for the development of habitat for openland wildlife, but if slopes are more than 15 percent, potential is poor. For development of habitat for woodland wildlife, these soils have good potential; for habitat for wetland wildlife, potential is very poor.

### 10. Savannah-Quitman

Nearly level to sloping, moderately well drained, loamy soils; some have a fragipan; on uplands and stream terraces

This map unit is in the eastern and southeastern parts of Rankin County. The landscape generally is nearly level to gently rolling but can include a few areas that are moderately steep. It is marked by broad ridges and nearly level, low terraces. The slope ranges from 0 to 8 percent.

This map unit makes up about 6.1 percent of the county. It is about 43 percent Savannah soils, 30 percent Quitman soils, and 27 percent soils of minor extent.

Savannah soils have a fragipan. They generally are in the raised, more sloping areas on uplands and stream terraces. These soils formed in loamy material. Quitman soils are in the low, smooth areas near drainageways, on uplands, and on stream terraces. These soils formed in loamy material. The minor soils in this map unit are Kirkville, Ora, and Tippah soils. Kirkville soils are moderately well drained and are on the flood plains. Ora and Tippah soils are moderately well drained and are on uplands.

Most of the acreage in this map unit is used as woodland or pasture. A small acreage is in crops.

In the gently sloping areas, Savannah and Quitman soils are well suited to row crops and small grains, and in the sloping areas, these soils are moderately suited to this use. The soils in this map unit are well suited to most grasses and legumes for hay and pasture.

Savannah soils are moderately suited to woodland. Windthrow and plant competition are moderate. Quitman soils are well suited to use as woodland. Wetness is a moderate limitation for equipment use.

Wetness is a moderate limitation to use of Savannah soils for most urban uses. Wetness and low strength as it affects local roads and streets are moderate limitation to use of Quitman soils for urban use.

Savannah and Quitman soils have good potential for the development of habitat for openland and woodland wildlife. In the nearly level areas, Quitman soils have a poor potential for development of habitat for wetland wildlife, and Savannah soils have a very poor potential.

### **Broad Land Use Consideration**

The soils in Rankin County vary widely in their suitabilities and limitations for major land uses. Kinds of soil limitations are indicated in general terms. The ratings of soil reflect the relative cost of practices to overcome the limitations and the hazard of continuing soil-related problems after practices are installed.

Kinds of land uses considered include cropland, pasture, woodland, urban development, and the development of habitat for wildlife. Cultivated farm crops grown extensively include cotton, soybeans, corn, and wheat. Woodland refers to land in trees. Urban areas include those used as residential, commercial, and industrial sites. Habitat for wildlife uses include habitat for openland wildlife, woodland wildlife, and wetland wildlife.

About 10 percent, or 49,853 acres, of Rankin County is used for cultivated crops, mostly soybeans, cotton, and wheat. Cropland is scattered throughout the county in areas of soils that are well suited to or moderately suited to row crops. These soils are mainly in map units 1, 3, 4, 5, 6, 8, and 10.

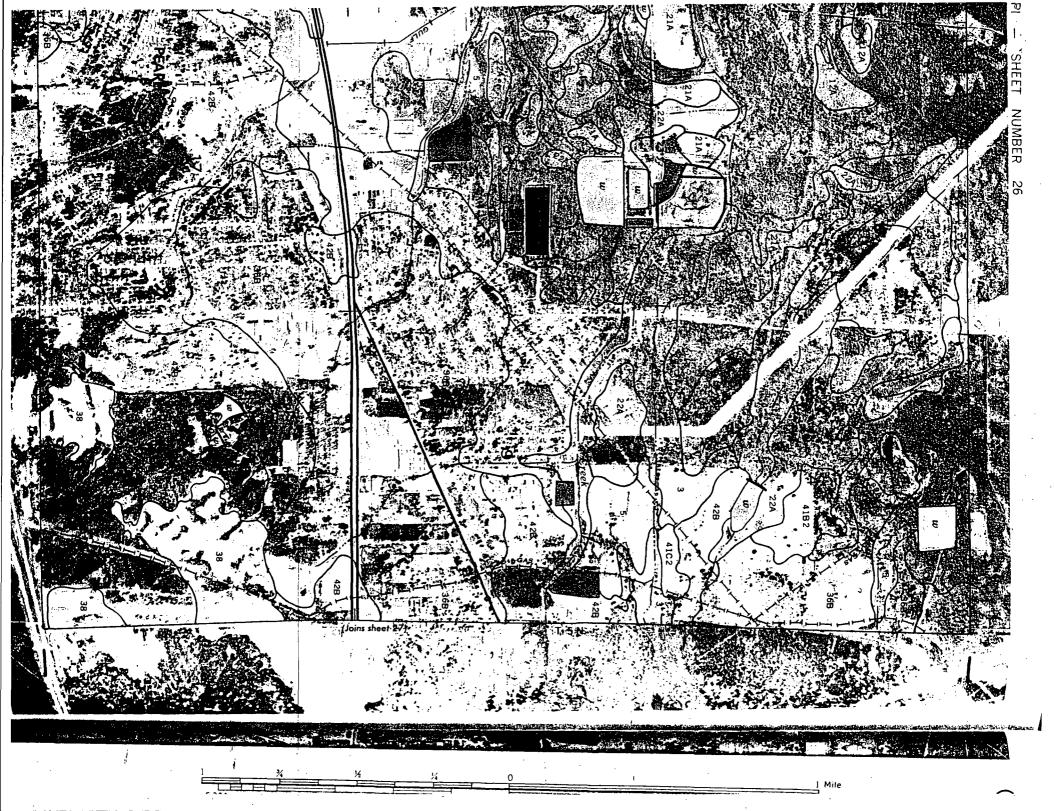
The soils in map units 3, 4, and 5 are occasionally flooded, mainly in winter and early in the spring. This flooding causes slight to moderate crop damage. The major soils in these map units are Urbo, Arkabutla, Oaklimeter, Gillsburg, Quitman, and Kirkville soils. Tippo and Guyton soils in map unit 1 are seasonally wet.

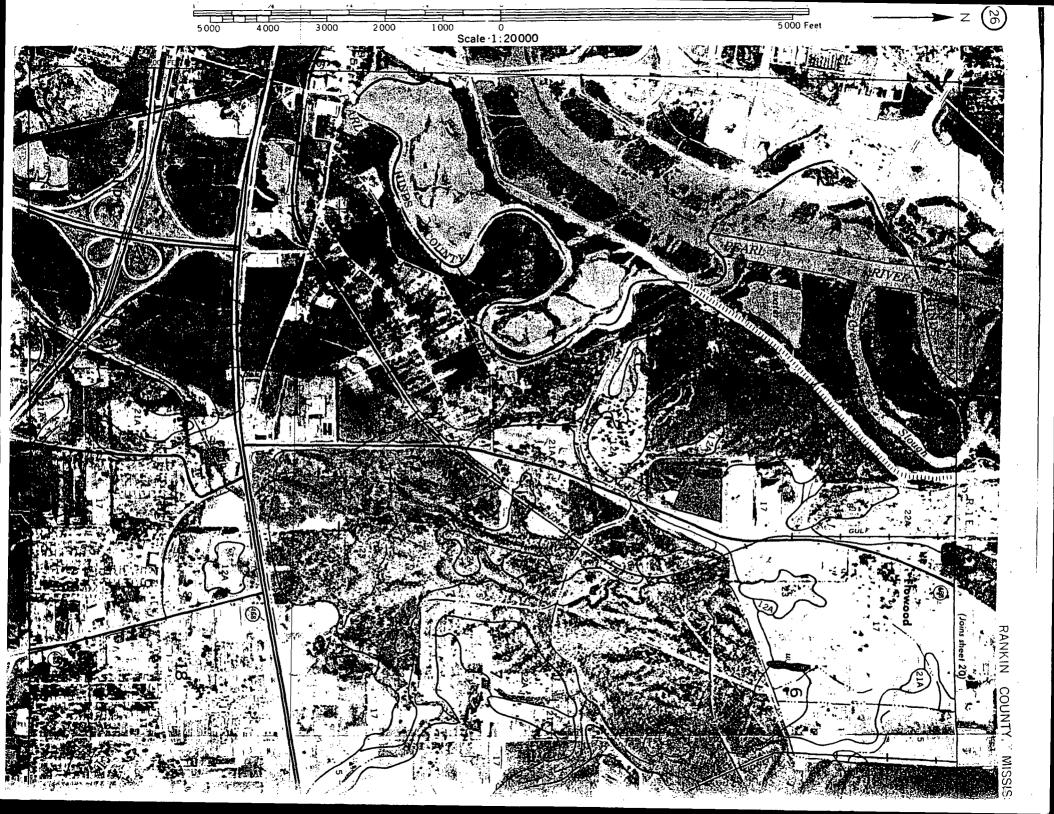
Erosion is a major hazard in growing crops on soils in map units 6, 8, and 10. Kipling, Falkner, Savannah, Providence, Tippah, and Quitman soils make up these map units. About 61 percent, or 310,000 acres, of the county is used as woodland. Soils in all map units are well suited to or moderately suited to trees. Some soils have a moderate to severe limitation for equipment use, but this limitation can be overcome by harvesting during the drier periods.

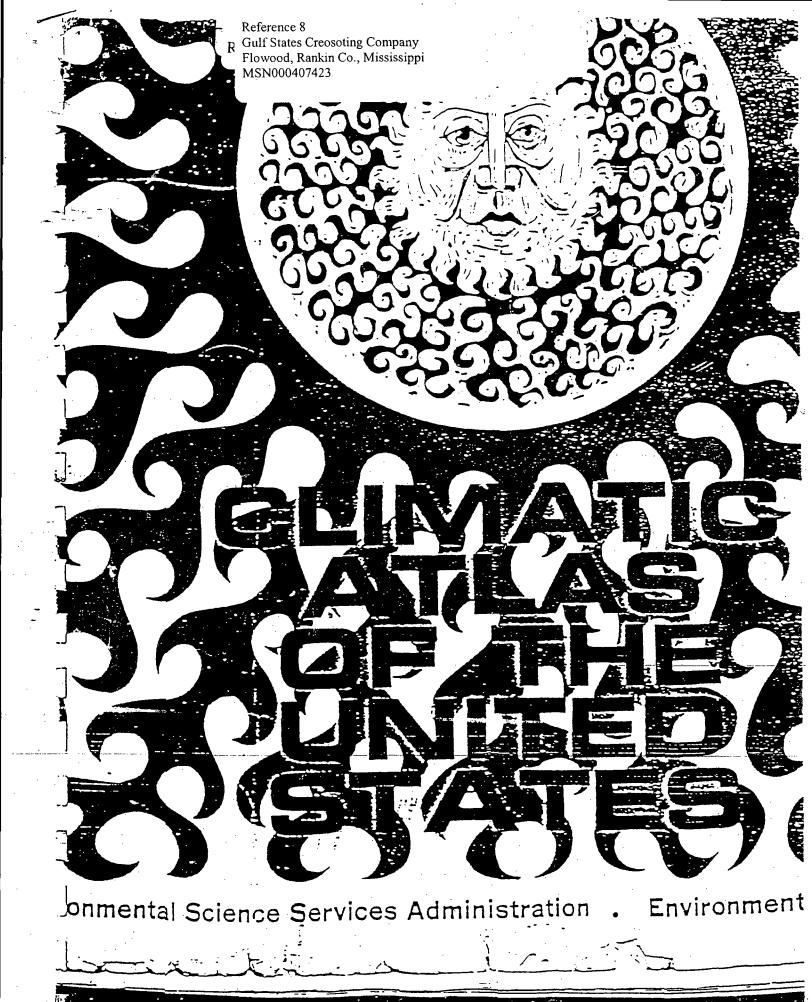
About 6 percent, or 33,176 acres, of the county is classified as urban or built-up land. Soils in map units 2, 3,4, and 5 that are on flood plains have severe limitations for urban use because of flooding. Quitman soils in map unit 5 are on higher elevations and are not subject to flooding. Soils in map unit 1 that are in protected areas have moderate limitations for urban use. Soils in map units 7 and 9 that are in hilly areas have severe limitations for urban use, mainly, because of the steepness of slope.

Soils in map units 8 and 10 have moderate limitations for urban use. High shrink-swell potential, low strength as it affects local roads and streets, and wetness are the main limitations of these soils for urban use. Most of the limitations can be overcome by special design and proper installation. The restricted permeability of Providence, Tippah, and Savannah soils is a limitation to use as septic tank absorption fields. This limitation can be partly overcome by enlarging septic tank absorption fields.

Kipling and Falkner soils in map unit 6 have severe limitations for urban use because of wetness and the high shrink-swell potential of the subsoil. Savannah soils have moderate limitations for urban use, mainly, because of wetness.









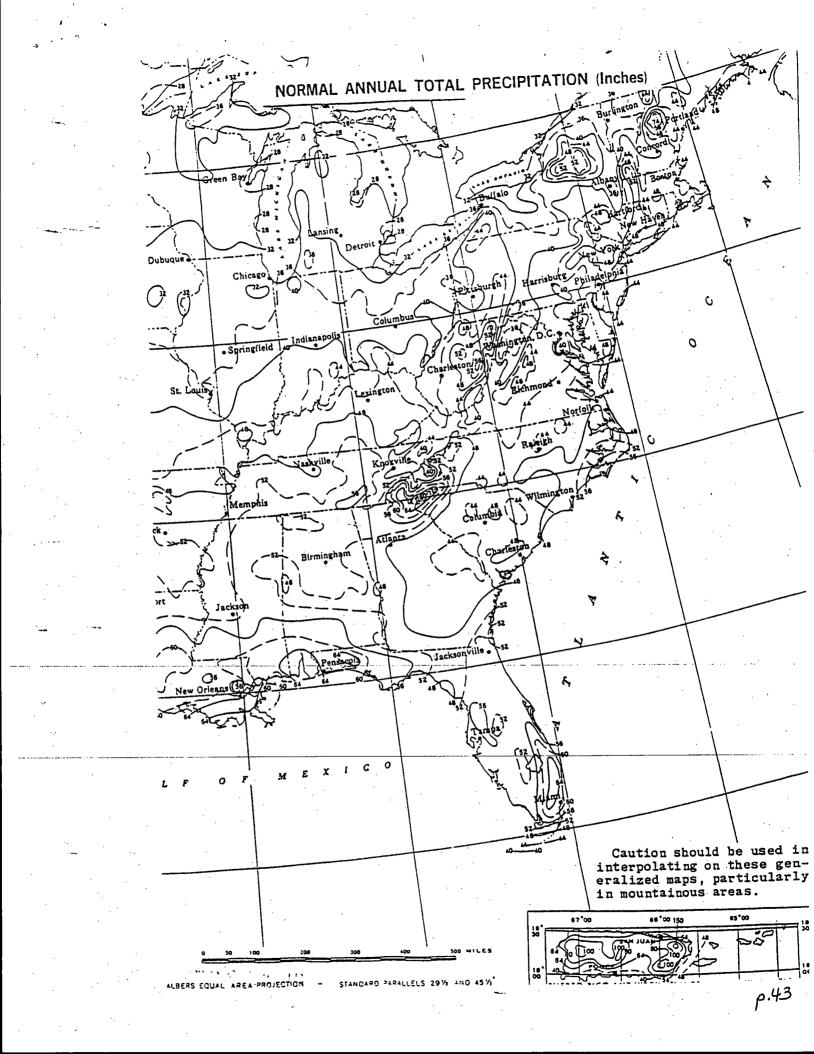
# U.S. DEPARTMENT OF COMMERCE C. R. Smith, Secretary

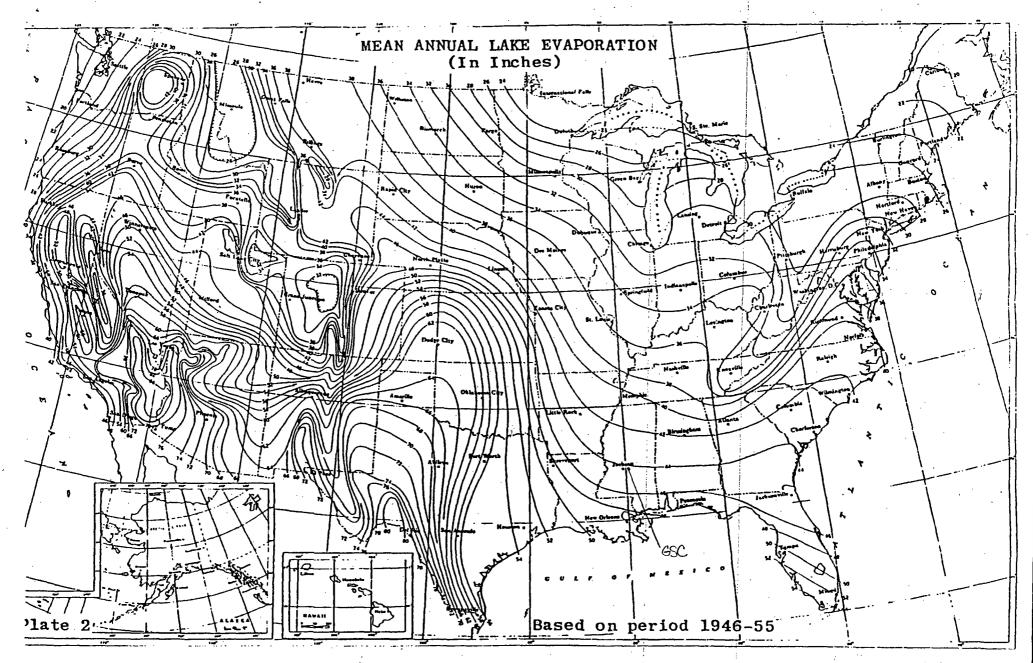
ENVIRONMENTAL SCIENCE SERVICES ADMINISTRATION Robert M. White, Administrator

ENVIRONMENTAL DATA SERVICE Woodrow C. Jacobs, Director

JUNE 1968

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U.S. DEPARTMENT OF COMMERCE LATHER II. HODGES, Sorretary

WEATHER BUREAU
F. W. REICHELDERFER, Chief

### TECHNICAL PAPER NO. 40

### RAINFALL FREQUENCY ATLAS OF THE UNITED STATES

### for Durations from 30 Minutes to 24 Hours and Return Periods from 1 to 100 Years

Prepared by
DAVID M. HERSHPIELD
Cooperative Studies Section, Hydrologic Services Division
for
Engineering Division, Soll Conservation Service
U.S. Department of Agriculture

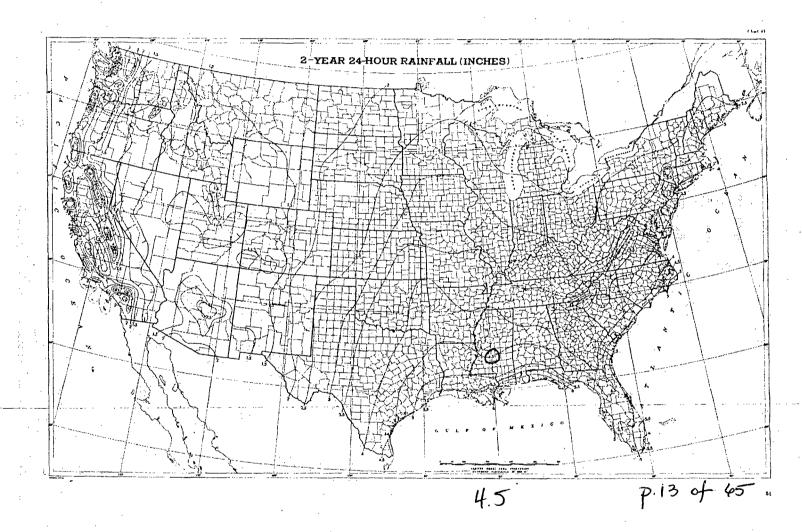


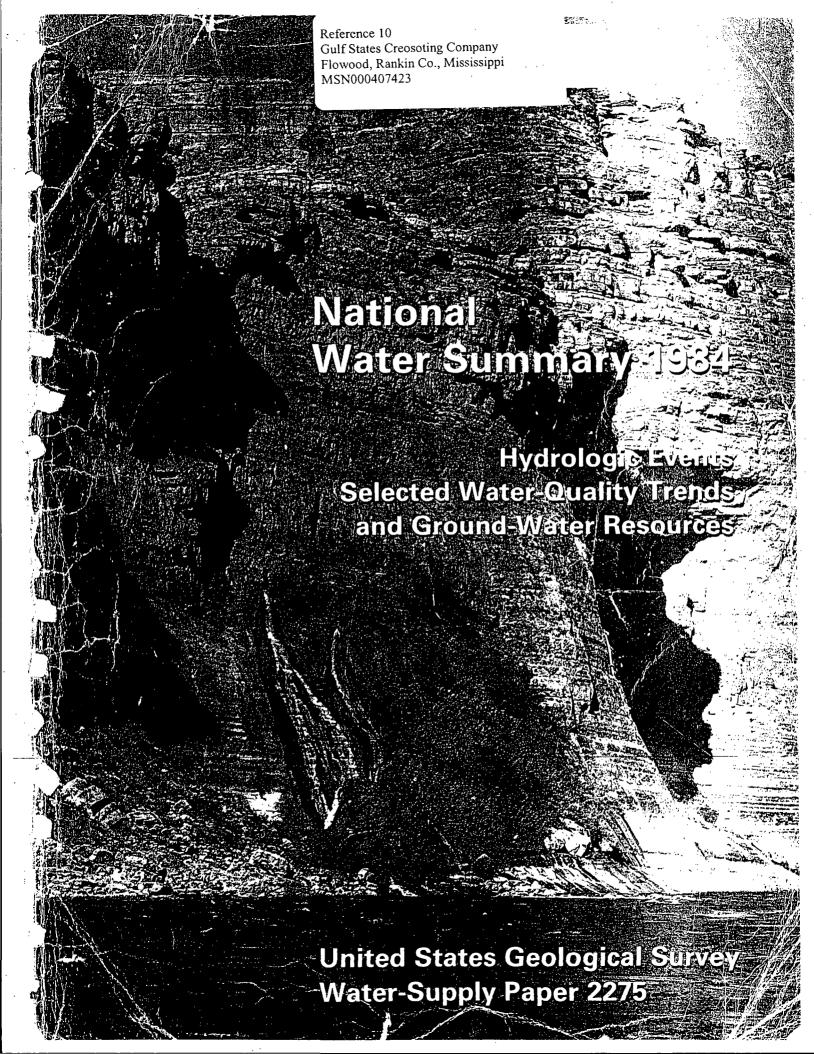
WASHINGTON, D.C.

May 1961

Repagnated and Repristed January 1963

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### Mississippi

### **Cround-Water Resources**

Ground water constitutes 54 percent of all freshwater used in Mississippi, serving the water supply needs of 93 percent of the population. The largest use of fresh ground water—54 percent of the total withdrawal—is for irrigation and aquaculture. Most of Jackson's public-water supply is withdrawn from the Pearl River but about 50 percent of the water used in the surrounding metropolitan area is from ground-water sources. Columbus and Meridian are converting from surface-water sources to wells. The nearly exclusive dependence on ground water for public-water supply is the result of statewide availability of aquifers that contain water of quality suitable for most uses and that are capable of supplying large yields [more than 300 gallons per minute (gal/min)] to wells. Ground-water withdrawals for various uses in 1980 and other related statistics are given in table 1.

### GENERAL SETTING

With the exception of an area of a few square miles in Tishomingo County, Mississippi lies entirely in the East Gulf Coastal Plain and is underlain by deposits of clay, sand, gravel, chalk, marl, and limestone. The oldest exposed strata are consolidated Paleozoic rocks that crop out only in a few valleys in Tishomingo County (fig. 1). Cretaceous strata in northern Mississippi dip and thicken southwestward. In central and southern Mississippi, the dip of the younger Eocene strata gradually becomes southward.

Much of the water that reaches the water table moves downdip westward to southwestward into the confined aquifers (fig. 1). Ground water moves westward into the northeastern Mississippi subsurface from Alabama. In southern Mississippi, some ground water flows into the subsurface of Louisiana or discharges into the Gulf of Mexico.

Precipitation in Mississippi is about 54 inches (in.) annually. Average monthly precipitation ranges from about 2.4 in. in October to about 6.2 in. in March. The late winter and spring rains provide an excess of water that results in high streamflow and periodic flooding. Infiltration from the Mississippi River and other streams reaches a maximum in the late spring.

About 50 percent of Mississippi's precipitation evaporates or is consumed by vegetation, about 40 percent runs off as streamflow, and about 10 percent infiltrates to the water table. Additional recharge of the ground-water reservoir is derived from infiltration of surface waters.

Several hundred gallons per minute can be obtained from wells completed in at least one aquifer nearly anywhere in the State. Throughout northwestern Mississippi and at places in the southern part of the State, well yields of several thousand gallons per minute are not unusual. Water-quality problems commonly are related to iron in solution and to acidic water. More troublesome in some areas, however, is the prevalence of color in ground water caused by the presence of organic

Table 1. Ground-water facts for Mississippi

[Withdrawal data rounded to two significant figures and may not add to totals because of independent rounding. Mgal/d = million gallons per day; gal/d = gallons per day. Source: Callahan, 1983]

Population served by ground water, 1980
Number (thousands) 2,33 Percentage of total population 9
ercentage of total population 9
rom public water-supply systems:
Number (thousands) 1,86
Number (thousands) 1,86 Percentage of total population 7
From rural self-supplied systems:
Number (thousands) 47
Percentage of total population 1
Freshwater withdrawals, 1980
Surface water and ground water, total (Mgal/d) 2,90
Ground water only (Mgal/d) 1.50
Ground water only (Mgal/d) 1,50 Percentage of total 5
Percentage of total excluding withdrawals for
thermoelectric power 8
Category of use
Public-supply withdrawals:
Ground water (Mgal/d) 23
Percentage of total ground water
Percentage of total public cumply
Percentage of total ground water 1 Percentage of total public supply 1 Per capita (gal/d) 12
Rural-supply withdrawals:
Domestic:
Ground water (Mgal/d) 2
Percentage of total ground water
Percentage of total rural domestic 10 Per capita (gal/d) 4
Per capita (gal/d) 4
Tivestock:
Ground water (Mgal/d) 8
Ground water (Mgal/d) 8 Percentage of total ground water 0
Percentage of total livestock
ndustrial self-supplied withdrawals:
Ground water (Mgal/d)43
Percentage of total ground water 2
Percentage of total industrial self-supplied:
Including withdrawals for thermoelectric power Excluding withdrawals for thermoelectric power
Excluding withdrawals for thermoelectric power
rrigation withdrawals:
Ground water (Mgal/d) 8
Percentage of total ground water
Percentage of total irrigation

Includes 264 Mgal/d for aquaculture use and 2.3 Mgal/d for waterfowl.

matter. Saltwater normally is present in the downdip parts of all aquifers; however, the base of freshwater extends to depths of more than 3,000 feet (ft) in some parts of the State (fig. 1). Saltwater intrusion has not been identified conclusively in coastal areas except locally where estuaries are connected hydraulically to shallow aquifers. Ground-water contamination from human activities is mostly restricted to oil-producing areas.

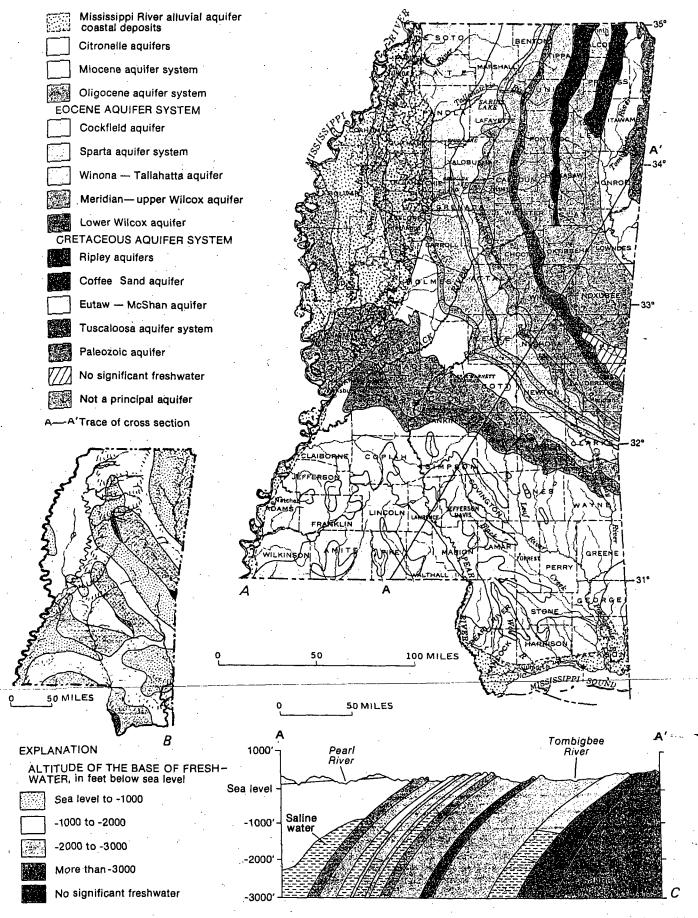


Figure 1. Principal aquifers of Mississippi. A, Geographic distribution. B, Altitude of the base of freshwater and physiographic diagram. C, Generalized cross section (A-A'). (See table 2 for a more detailed description of the aquifers. Sources: A, Modified from Bicker, 1969. B, Gandi, 1982; Raisz, 1954. C, Compiled by E. H. Boswell from U.S. Geological Survey files.)

### 270 National Water Summary—Ground-Water Resources

Table 2. Aquifer and well characteristics in Mississippi

[Gal/min = gallons per minute; mg/L = milligram per liter; ft = feet. Sources: Reports of the U.S. Geological Survey, Mississippi Bureau of Land and Water Resources, and Mississippi Research and Development Center]

quifer name and description Depth (ft) Yield (gal/min)				Remarks	
quita manta ana accomption	Common range	May exceed	Common range	May exceed	- Tomains
Mississippi River alluvial aquifer: Sand, gravel, silt, and clay. Semiconfined.	50 - 140	200	500 - 3,000	5,000	Water hard, iron in solution generally exceeds 1.0 mg/L. Susceptible to pollution. Source of public water supply at Vicksburg (location 15, fig. 2).
Citronelle aquifers: Sand, gravel, silt, and clay. Generally unconfined.	50 - 200	250	50 – 300	500	Water soft, acidic, iron in solution generally exceeds 0.3 mg/L. Dissolved solids concentrations generally lower than 100 mg/L. Source for several public water supplies in southern part of State. Susceptible to pollution. Equivalent to Pliocene-Miocene aquifer in Alabama, Pleistocene aquifer in Louisiana.
Miocene aquifer system: Sand, clay, gravel, and silt. Generally confined.	50 - 1,500	2,400	50 – 1,500	5,000	Includes Graham Ferry, Pascagoula, and Hattiesburg Formations and Catahoula Sandstone. Water soft, sodium bicarbonate type; locally, iron exceeds 0.3 mg/L. Contaminated by oilfield brine locally. Principal source for public water supplies in southern one-third of State. Equivalent to Pliocene-Miocene aquifer in Alabama and Louisiana.
Oligocene aquifer system: Limestone, sand, silt, and clay. Generally confined.	150 - 1,000	1,200	10 - 150	400	Includes Vicksburg Group and Forest Hill Sand. Water soft, slightly alkaline. Source for a few public water supplies in south-central part of State. Part of Oligocene-Eocene aquifer in Alabama. Confining unit in Louisiana.
Eocene aquifer system:  Cockfield aquifer: Sand, silt,  clay, and lignite. Generally  confined.	100 - 1,000	1,200	10 - 1,000	1,500	Water hard near outcrop, sodium bicarbonate type elsewhere. Locally, iron concentration exceeds 0.3 mg/L and color is more than 20 units. Largest withdrawal is for public water supply at Greenville (location 16, fig. 2). Part of Tertiary sand aquifer in Tennessee, Oligocene-Eocene aquifer in Alabama.
Sparta aquifer system: Sand, silt, clay, and lignite. Generally confined.	100 - 1,500	2,000	10 – 1,000	3,000	Water soft, sodium bicarbonate type. Locally iron concentration exceeds 0.3 mg/L and color is more than 20 units. Contaminated by oil-field brine locally. Source for many public water supplies in central and northwestern Mississippi. Part of Tertiary sand aquifer in
					Tennessee, Oligocene-Eocene aquifer in Alabama.
Winona-Tallahatta aquifer: Glauconitic sand and clay. Generally confined.	100 - 1,000	1,200	10 - 400	500	Water soft. Locally, iron concentration exceeds 3.0 mg/L, and color is more that 20 units. Source for public water supply for several small municipalities. Part of Tertiary aquifer in Tennessee. Oligocene-Eocene aquifer in Alabama. Confining unit in Louisiana.
Meridian-upper Wilcox aquifer: Sand, silt, clay, and lignite. Generally confined.	100 - 1,800	2,000	100 - 2,000	2,500	Water soft, acidic in the north. Locally iron concentration exceeds 0.3 mg/L, and color is more than 20 units. Source for many public water supplies in centra and northwestern Mississippi. Largest withdrawal is at Greenwood (location 16 fig. 2). Part of Tertiary sand aquifer in Tennessee, Oligocene-Eocene aquifer in Alabama, and Wilcox-Carrizo aquifer in Louisiana.

Table 2. Aquifer and well characteristics in Mississippi-Continued

	Well characteristics				
Aquifer name and description	Depti	h (ft)	Yleld (g	al/min)	Remarks
	·	Common range	May exceed	Common range	May exceed
Lower Wilcox aquifer: Sand, silt, clay, and lignite. Generally confined.	100 – 2,100	3,000	100 ~ 1,500	2,000	Water soft. Locally, iron concentration exceeds 0.3 mg/L. Contaminated by oil-field brine locally. Source for public water supplies throughout central and northwestern Mississippi. Largest
		,			withdrawal is at Meridian (location 22, fig. 2). Equivalent to Nanafalia-Clayton aquifer in Alabama and part of Tertiary sand in Tennessee.
cretaceous aquifer system: Ripley aquifers: Sand, clay, sandstone, and limestone. Generally confined.	50 - 1,100	1,800	10 - 300	400	Water hard near outcrop, soft at depth. Source for several small public water supplies in extreme northern part of State. Part of Cretaceous aquifer in Tennessee.
Coffee Sand aquifer: Sand, clay and sandstone. Generally confined.	50 - 1,000	2,000	10 – 400	.500	Water hard near outcrop, soft at depth. Source for several small public water supplies in extreme northern part of State. Part of Cretaceous aquifer in Tennessee.
Eutaw-McShan aquifer: Sand and clay. Generally confined.	100 - 1,500	1,800	10 - 500	600	Water hard near outcrop, soft at depth. Locally fluoride exceeds 1.0 mg/L. Source for numerous public water supplies in northern part of State. Largest withdrawals are at Tupelo and in Monroe County (locations 9 and 23, fig. 2). Equivalent to Eutaw aquifer in Alabama and part of Cretaceous aquife in Tennessee.
Tuscaloosa aquifer system: Sand, gravel, silt, and clay. Generally confined.	100 – 2,000	2,400	50 – 1,500	2,000	Includes Gordo and Coker Formations, and locally, beds of Early Cretaceous age. Water soft to slightly hard, small dissolved-solids concentrations.
					Locally iron exceeds 0.3 mg/L. Source for numerous public water supplies in northwestern Mississippi. Largest withdrawals are in Columbus area and Monroe County (location 23, fig. 2). Equivalent to Tuscaloosa aquifer in Alabama.
Paleozoic aquifer: Limestone, chert, and clay. Generally confined.	100 - 600	1,000	100 - 900	1,000	

The southwestward dip of the strata and the overlap of freshwater in successively younger aquifers southward result in the availability of two or more separate aquifers for development in most places (fig. 1). Examples are use of both the Tuscaloosa aquifer system and the Eutaw-McShan aquifer at localities in the northeast; the Cockfield, Sparta, and Meridian-upper Wilcox aquifers in some mid-State localities; and the Meridian-upper Wilcox and lower Wilcox aquifers in many areas. Some geologic formations include two or more extensive water-bearing zones that function as a single system when considered on a regional basis (Sparta aquifer system). Other water-bearing formations are directly connected hydraulically and function as a single aquifer (Eutaw-McShan aquifer).

### PRINCIPAL AQUIFÉRS

Except for the chert aquifer of Paleozoic age, which is the source of water for several public-water supplies in Alcorn and Tishomingo Counties, all principal aquifers in Mississippi consist of unconsolidated sand or sand and gravel strata that are irregular in thickness and physical character and exhibit extreme variation in their capability to store and transmit water (Wasson, 1980). The principal aquifers are discussed below and in table 2; their areal distribution is shown in figure 1.

### MISSISSIPPI RIVER ALLUVIAL AQUIFER

The extensive Mississippi River alluvial aquifer in the Delta area of northwestern Mississippi is an extremely prolific

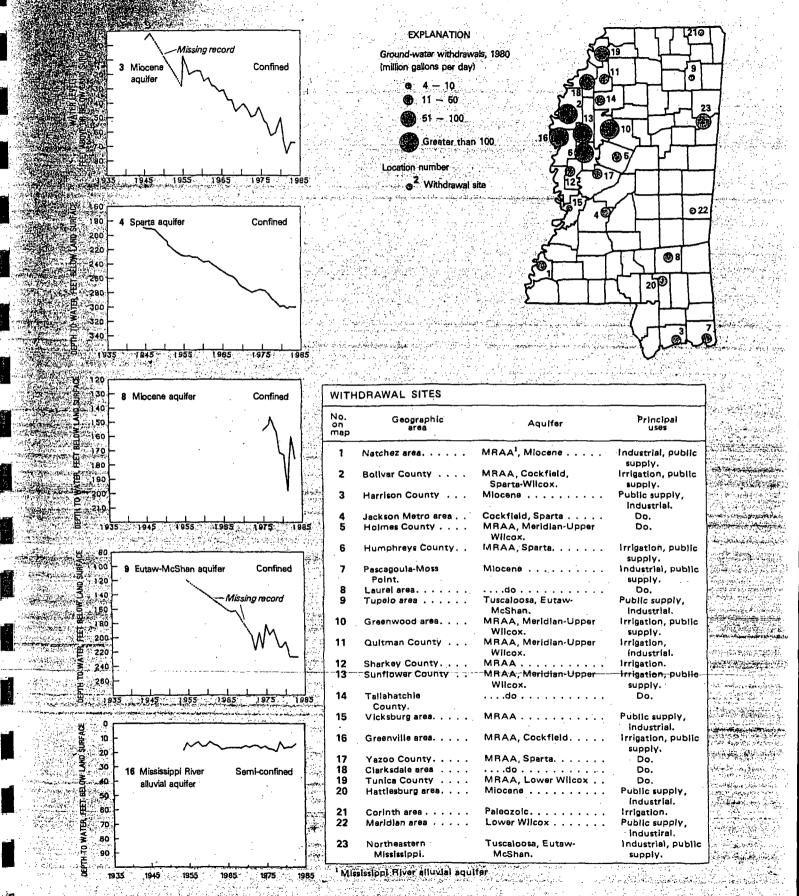


Figure 2. Areal distribution of major ground-water withdrawals and graphs of annual greatest depth to water in selected wells in Mississippi. (Sources: Withdrawal data from Callahan, 1983; water-level data from U.S. Geological Survey files.)

source of water that is used for irrigation, aquaculture (principally catfish farming), industrial cooling, and for one public supply (Vicksburg). The alluvium averages about 140 ft in thickness. Generally, the uppermost 20 to 30 ft is clay or other fine-grained material; underlying sand and gravel beds form the aquifer. Wells about 120 ft deep that produce 2,000 to 3,000 gal/min can be constructed nearly anywhere in the Delta.

The alluvial aquifer is recharged by the Mississippi River and smaller streams and, to a lesser extent, by direct infiltration of precipitation (Sumner and Wasson, 1984). Recharge also occurs on the east side of the delta where streams enter from the Bluff Hills and where water-bearing zones in the deeper aquifers are in contact with the alluvium.

### CITRONELLE AQUIFERS

The Citronelle aquifers overlie older aquifers in southern Mississippi and are used for some public and industrial wells and extensively for small domestic and farm wells. The Citronelle originated as an extensive surficial fluvial deposit that has now been greatly dissected by streams. The relatively flat-lying, very permeable beds are a source of water for springs and seeps that sustain the low flow of streams and transmit recharge to underlying confined aquifer subcrops.

### MIOCENE AND OLIGOCENE AQUIFER SYSTEMS

The aquifers in southern Mississippi partly underlie the Citronelle aquifers and are separated from the underlying Eocene aquifers by several hundred feet of clay. In descending order, these aquifers are present in the Graham Ferry Formation, the Pascagoula Formation, the Hattiesburg Formation, the Catahoula Sandstone, the Vicksburg Group, and the Forest Hill Sand. All except the Vicksburg and the Forest Hill aquifers, which form the Oligocene aquifer system, are included in the Miocene aquifer system.

Some water wells in the Miocene aquifers are about 2,000 ft deep and the deepest well reaches 2,400 ft; however, geophysical logs made of oil tests in Hancock County show that freshwater extends to slightly more than 3,000 ft below sea level (fig. 1). Water wells about 1,000 ft deep on some of the barrier islands that form Mississippi Sound confirm that freshwater aquifers extend gulfward beyond the shoreline (Brown and others, 1944).

### **EOCENE AQUIFER SYSTEM**

The Eocene aquifers, exposed at the surface in north-central, northwestern, and central Mississippi, extend in the subsurface to the west, southwest, and south, and contain freshwater in about 50 percent of the State (Wasson, 1980). Included are the Cockfield, the Winona-Tallahatta, the Meridian-upper Wilcox, and the lower Wilcox aquifers and the Sparta aquifer system. All are regional in extent, and all except the Cockfield and lower Wilcox merge northward into a single aquifer south of Memphis, Tennesseee. The deepest water well in Mississippi (2,760 ft) taps the fower Wilcox aquifer in northern Wayne County. Geophysical logs made in oil test wells show that freshwater in this aquifer extends more than 3,000 ft below sea level in Smith County (fig. 1).

#### CRETACEOUS AQUIFER SYSTEM

The Cretaceous aquifers contain freshwater in about one-fourth of the State (Boswell, 1963). The outcrop area is in northeastern Mississippi. Cretaceous aquifers include the Ripley, the Coffee Sand, and the Eutaw-McShan aquifers, and the Tuscaloosa aquifer system. The Eutaw-McShan and Tuscaloosa aquifers extend into Alabama. The Ripley and the Coffee Sand aquifers, which are restricted to northern Mississippi generally north and west of Tupelo, extend into Tennessee and Arkansas. Freshwater extends to depths that exceed 3,000 ft below sea level in some areas more than 80 miles from the recharge areas (fig. 1), and some water wells exceed 2,000 ft in depth. The deepest wells that tap the Tuscaloosa aquifer system are located in the outcrop area of the Eocene aquifers (fig. 1).

#### PALEOZOIC AQUIFER

The Paleozoic aquifer consists of the upper part of weathered, faulted limestone and chert; the aquifer is overlain by Cretaceous deposits in extreme northeastern Mississippi. Present development of the aquifer is restricted to Alcorn and Tishomingo Counties where well depths range from 100 to 600 ft (Wasson, 1980). Wells produce as much as 1,000 gal/min where large declines in water levels have not occurred. The water is moderately hard and, at some sites, contains more than 0.3 milligram per liter (mg/L) of iron.

### GROUND-WATER WITHDRAWALS AND WATER-LEVEL TRENDS

Pumping centers that produce 4 million gallons per day (Mgal/d) or more of ground water are shown in figure 2. About three-fourths (1,143 Mgal/d in 1980) of the ground water used in Mississippi is pumped in the northwestern part of the State (fig. 2) from the Mississippi River alluvial aquifer for irrigation and aquaculture (Callahan, 1983). Water levels in the aquifer fluctuate seasonally, reaching high levels in the spring after recharge and declining to the lowest point in the fall following irrigation withdrawals and normal seasonal decline. The hydrograph for the alluvial aquifer near Greenville (location 16, fig. 2), indicates recovery of water levels in the aquifer after 1975; however, in some other areas (locations 2, 10, and 12), water levels lowered by the combination of drought conditions and unprecedented use of water for irrigation and aquaculture have not recovered in some areas (Sumner and Wasson, 1984).

Confined aquifers in Mississippi generally have shown a regional decline of about 2 ft annually during the last 30 years. The declines have attracted attention where pumping is concentrated in aquifers that are only a few hundred feet deep and pumping occurs near the top of the aquifer (locations 7, 9, and 21, fig. 2). The effects of water-level declines have elicited less concern in other areas where several hundred feet of available drawdown remains. The most pronounced water-level declines have been in the Paleozoic aquifer at Corinth (location 21), the Eutaw-McShan aquifer at Tupelo (location 9), the Sparta aquifer system at Jackson (location 4), and the Miocene aquifer system at Natchez, Pascagoula, Laurel, and Hattiesburg (locations 1, 7, 8, 20). Water-level recovery

during the last several years at locations 7 and 8 is due to reductions in withdrawal and changes in pumping distribution.

### **GROUND-WATER MANAGEMENT**

The 1956 omnibus water law passed by the Mississippi Legislature specifically excluded subsurface waters. It was not until 1976 that a ground-water bill, codified now as Sections 51-4-1 et. seq., Mississippi Code Annotated, 1972 (James I. Palmer, Jr., Governor's Office of Economic Development and Natural Resources, written commun., 1984) was enacted. The concept of "capacity use areas," wherein well spacing, well depths, and withdrawal rates are regulated, is the mechanism provided for dealing with areas having identifiable ground-water-supply problems. The major limitations of the 1976 Act are that it addresses only withdrawals in excess of 50,000 gallons per day (gal/d) and excludes agricultural and oil and gas uses. In 1983, the State legislature created the Mississippi Water Management Council to reexamine completely all State laws pertaining to surface and subsurface waters and to report recommended amendments to the 1985 session.

The Mississippi Department of Natural Resources administers and enforces, through its Bureau of Land and Water Resources, not only the 1956 surface-water and 1976 ground-water statutes but also the 1966 Water Well Drillers Licensing Act. Primacy in permitting waste injection in Mississippi (other than in connection with oil and gas production) has been assigned to the Department's Bureau of Pollution Control, which also has responsibility for permitting and monitoring hazardous-waste sites. On June 27, 1984, Mississippi became the second State to be given final authorization to operate its own hazardous-waste program. Primacy for permitting oil field waste injection has not been delegated by the U.S. Environmental Protection Agency (as of December 1984).

The Department's Bureau of Geology, basically a research organization, is authorized to investigate and report on water resources. The Mississippi State Board of Health ensures that public-water supplies meet chemical, bacteriological, and other standards.

Water-resources investigations in Mississippi are conducted cooperatively by the U.S. Geological Survey with the Mississippi Department of Natural Resources, 10 other State and local agencies and municipalities, and five Federal agencies.

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## Rankin County Geology and Mineral Resources

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**BULLETIN 115** 

## MISSISSIPPI GEOLOGICAL, ECONOMIC AND TOPOGRAPHICAL SURVEY

WILLIAM HALSELL MOORE
Director and State Geologist

JACKSON, MISSISSIPPI
1971

PRICE \$3.00

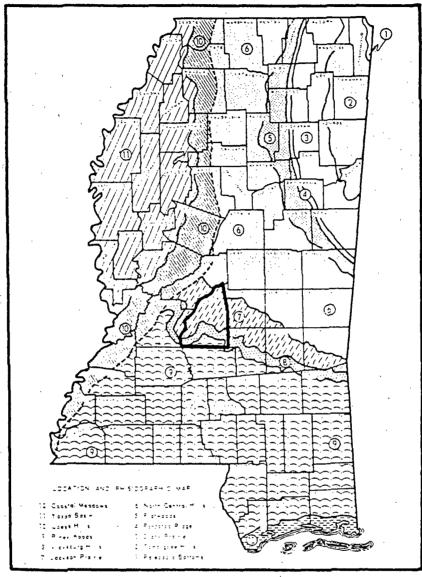


Figure 3.—Physiographic provinces of Mississippi. After R. R. Priddy.

Lowe<sup>21</sup> included the whole southern half of Mississippi south of the Jackson Prairie belt in a unit he designated the Long Leaf Pine Hills. This unit is now commonly known as the Southern Pine Hills. Priddy<sup>22</sup> further divided this unit into 3 dis-

The Paleocene-Eocene boundary has not been satisfactorily established in the subsurface of Mississippi. A time-rock unit concept, based on restricted faunal zones, would greatly expand the thickness of the Paleocene beds to include substantial amounts of the overlying strata now classified as Wilcox (Eocene). Because of its general acceptance as a rock unit, perpetrated mainly by the need for a reliable and easily identifiable electrical log correlation, the Midway is treated as a rock unit. In this sense the Paleocene sediments are restricted to a usal calcareous unit, the Clayton formation and an upper shale unit, the Midway shale.

### Wilcox Group

The Wilcox group in Rankin County consists of an undifferentiated mass of complexly interstratified clays, clay shales, sandstones, silts and a few thin marls. The Wilcox varies from 1100 feet to approximately 1300 feet in thickness over the Jackson Dome. Off the dome, Wilcox sediments attain a thickness of 2830 feet in the more southern part of the County.

Wilcox sandstones are very fine- to coarse-grained. Quartz pebbles are commonly associated with the coarser sandstones. Most of the sandstones contain at least some lignitic or carbonaceous material and are also commonly slightly micaceous. The sandstones and silts are white, pale-gray, light-gray and light-green, the varying amounts of clay, silt and carbonaceous matrix materials producing the various colors. Some of the sandstones are glauconitic and calcareous. In some cases, these glauconitic sandstones were associated with the thin pale-gray to light-gray sandy marls.

The clays and clay shales are pale-gray, light-gray, gray and dark-gray. The gray and dark-gray clays and clay shales generally contain finely divided carbonaceous and lignitic material. Numerous seams of lignite were noted.

Any Meridian sand equivalents in the subsurface of Rankin County are included in the Wilcox group rather than in the Claiborne group. The inclusion of the Meridian sand in the Wilcox group is common practice when it is not easily separable from the subjacent Wilcox sediments.

#### Winona Formation

The Winona consists of a sequence of interbedded pale-gray and pale-grayish-white silty glauconitic chalks, sandy marls and minor amounts of light-gray and light greenish-gray slightly calcareous and fossiliferous clay shales and clays.

The Winona formation is only 10 to 15 feet thick over the Jackson Dome. Thickening occurs in all directions off the dome and up to 65 feet of Winona, restricted as it is in this report, have been recorded in the eastern parts of the County.

#### Zilpha Formation

The Zilpha varies from a minimum of 200 feet in thickness on the Jackson Dome to a maximum of 420 feet in the southwestern part of the County. Facies changes produced by local depositional environments, as well as structural thinning over the Jackson Dome account for the divergence in formation thickness.

The lower part of the Zilpha formation consists of a sequence of gray slightly fossiliferous clay shales often containing finely disseminated lignite or carbonaceous material. The basal strata of the Zilpha is slightly glauconitic and calcareous at the contact with the glauconitic chalks and marls of the underlying Winona formation.

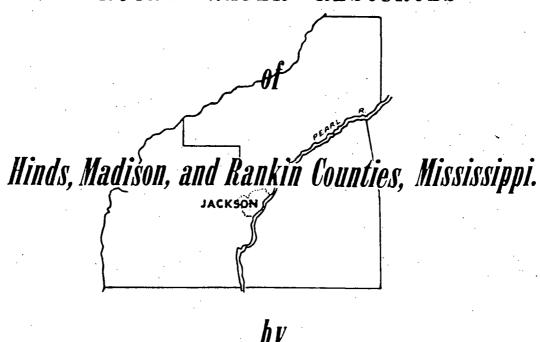
The upper part of the formation is variably sandy and the clay shales tend to be silty and slightly micaceous. A fluctuating succession of one facies of deposition by another resulted in an interstratified sequence of clay shales and very fine- to medium-grained sands. These sands are finer than the basal sands of the overlying Kosciusko formation. Some of the finer sandstones in several wells were slightly glauconitic. Traces of reddish-brown slightly sandy and glauconitic clay ironstones were also-noted in several wells. The few available sets of cuttings through this interval were of such poor quality that any distribution pattern of the glauconitic sandstones and clay ironstones could not be determined.

The top of the Zilpha is placed at the base of the lowest massive sand sequence of the Kosciusko formation. While the upper part of the Zilpha may be quite sandy, these sands are

Reference 12 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423

STATE OF MISSISSIPPI BOARD OF WATER COMMISSIONERS BULLETIN 64 - 1

## GROUND - WATER RESOURCES



E. J. Harvey, J. A. Callahan, and B. E. Wasson

Prepared by the United States Geological Survey
In cooperation with the
City of Jackson and the Mississippi Board of Water Commissioners
April 1964

Ву

E. J. Harvey, J. A. Callahan, and B. E. Wasson

#### ABSTRACT

The Jackson, Mississippi, metropolitan area, consisting of Hinds, Madison, and Rankin Counties, possesses abundant reserves of ground water of good quality. This area, comprising 2,400 square miles, is on the eastern flank of the Mississippi embedient. Regional strike of the formations is northwest and dip is southwest. The Jackson dome is the main structural feature, and formation strike and dip in the vicinity of the dome are at variance with regional trends.

Fresh ground water occurs to depths greater than 2,000 feet almost everywhere in the area. The major sources of water supply, potentially and present, are the Wilcox Group, Sparta Sand, Cockfield Formation, Forest Hill Sand, and Catahoula Sand stone. Small supplies of hard water can be obtained from surficial deposits underlying stream terraces and the flood plains of the Big Black and Pearl Rivers.

The Wilcox Group contains a large reserve of soft water that has been tapped by only a few small-supply wells. The water is more highly mineralized and is warmer than that in the shallower aquifers. It is of good quality in Madison and northern Rankin Counties, but the quality deteriorates down the dip in Hinds County. Treatment to reduce color may be required for Wilcox water supplies.

The Sparta Sand is the principal aquifer from the standpoint of total pumpage. Uplift associated with the Jackson dome centered beneath the city of Jackson has raised the Sparta about 600 feet and made it readily accessible to development. Sparta wells 600 to 1,000 feet deep in Jackson have yields ranging from 200 to more than 1,000 gpm (gallons per minute). About 10 mgd (million gallons per day), 40 percent of the water used in Jackson, is pumped from the Sparta. It is probable tha additional properly designed well fields can withdraw as much as 5 mgd each without lowering the water level below feasible pumping depths.

The Cockfield is the most widely tapped water-bearing formation, chiefly because it is the shallowest aquifer in much of the area. Municipal water supplies for several small towns are obtained from the Cockfield, some wells yielding as much as 500 gpm.

Water from the Sparta Sand and Cockfield Formation is a soft, sodium bicarbonate type having a dissolved-solids content generally less than 400 ppm (parts per million Iron is a problem in some places, especially in the Cockfield water, and color is mor commonly objectionable in that aquifer.

The Forest Hill Sand and Catahoula Sandstone supply three small municipal water supplies and many domestic wells in the southern part of the area. Dissolved solids average about 500 ppm in water from these aquifers. Water from the Forest Hill generally is colored, and silica is high in the few Catahoula water samples analyzed.

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.- 6 -- 35 -- 36 dome. The aquifers described in this report range in geologic age from early Ecc to Recent. Monroe (1954, p. 50-110) gave detailed descriptions of the outcroppir rocks in Rankin County and in the vicinity of the Jackson dome in Hinds and Madis Counties. The outcropping rocks of Madison County were described by Priddy (1960)

Owing to the regional southwestward dip, the older beds crop out in the nort eastern part of the area and dip beneath the surface. Progressively younger form tions are at the surface in belts a fraction of a mile to as much as 30 miles wid depending on formation thickness and dip. The belts of outcrop (fig. 1) strike northwest, except in the vicinity of Jackson where they form incomplete concentri rings on the flanks of the Jackson dome. Beds that reappear at the surface on the crest of the dome would be 600 to 700 feet below the surface were the dome nonexi

The oldest units described in this report are sands of the Wilcox Group, as are the oldest strata containing fresh water in these counties. In ascending ord the chief aquifers above the Wilcox Group are the Sparta Sand, Cockfield Formatic Forest Hill Sand, and Catahoula Sandstone. Fresh water occurs generally to depth much as 2,000 feet and possibly to a considerably greater depth in some localitie Lithologic and hydrologic characteristics of the various rock units underlying th area are summarized in table 1.

In the geologic column the water-bearing sand beds are interbedded with shal both marine and continental origin, fossiliferous sandy limestone, and calcareous sandstone. Strata deposited under generally marine conditions consist mostly of and they form aquicludes. The aquicludes are widespread and more uniform in thic than the aquifers. Sand beds composing the aquifers were deposited chiefly in a taic environment, and they pinch out or lens with clay to such an extent that san make up only 5 or 10 percent of the entire thickness of the formation in one loca and more than 50 percent in another. Lignitic material is common in several plac However, as a unit an aquifer such as the Sparta Sand, which contains several wat bearing beds, can be traced across the entire area.

#### STRUCTURE

Jackson is on the eastern flank of the broad Mississippi embayment syncline, part of the Gulf Coastal Plain province. The Tertiary formations dip southwest f the outcrop area in the northern and central parts of the State toward the embaym axis, which lies immediately west of Hinds County and generally parallels the Mis sippi River. West of the axis, the formations rise to the outcrop areas in Arkan and Louisiana. Dips in the project area usually are 15 to 25 feet per mile. Loc dips are greater and they vary from the normal southwesterly direction. The regidip of the Tertiary rocks is interrupted by the Jackson dome, where the strata ar arched upward several hundred feet. The strike of the formations is northwest ex in the vicinity of the dome.

The sections, contour maps, and isopachous maps in the report demonstrate th pronounced effect of the Jackson dome on the attitude, thickness, and depth of aq fers. The distribution of detrital materials on the flanks of the dome has a con siderable effect on the availability of water supplies. There is a probability t slow, continuous movement and warping took place during deposition as loading by mentation increased around the dome. Because of the raised position of the dome, a part of the detrital materials available to the area was deposited on the struc

#### er Quality

Six of the seven wells producing water from the Wilcox Group are on the crest of a Jackson dome or high on its flanks. There is considerable variation in the quaty of water from these wells. The water is in general a soft, colored, mixed sodium carbonate-sodium chloride type (table 4).

Electric logs indicate that fresh water is available in all of Madison County, a northeastern part of Hinds County, and the northern half of Rankin County. The per sand in the Wilcox contains fresh water over a larger part of the area than do a deeper sands. In the northern part of Madison County fresh water is available a depth of 2,200 feet, or to within 500 feet of the base of the Wilcox. On the me and its flanks in the vicinity of Jackson, the base of fresh water is about 1,800 at deep. South of the Jackson dome the Wilcox reaches depths at which all the water salty. A line corresponding with the -1,800-foot contour (fig. 7) may be consided the limit of fresh water southwest of Jackson.

#### Development of Water Supplies

The potential of the Wilcox as a source of water supply in this area can only be eorized. In counties to the northeast, the results of several pumping tests indite that sands in the Wilcox propably are as permeable as those in the Sparta.

#### Tallahatta Formation, Winona Sand, and Zilpha Clay

The Tallahatta Formation, whoma Sand, and Zilpha Clay are mainly of marine igin. The sediments constant in ascending order, of: light-gray to greenish-gray ugh clay and siltstone of the Tallahatta, glauconitic sand and sandstone of the nona, and reddish-brown shay and shale of the Zilpha.

These units are essentially non-water-bearing and constitute an aquiclude (relavely imperments bed) in the area. They are not differentiated in this report. A tailed lithologic and stratigraphic sequence of the beds was given in the basic-ta report (Harvey, Calla an, and Wasson, 1961, p. 60-61).

The three formations aderlie the entire area at depths that range from 650 feet northeastern Madison County to 2,600 feet in southwestern Hinds County. The aggrete thickness is about 300 feet over the Jackson dome. Thickness in other parts of a area ranges from 420 to 570 feet.

#### Sparta Sand

#### Distribution and Character

The Spart sand, formerly known as the Kosciusko Formation (Thomas, 1942, p. 40), in the subsurface throughout the three counties (figs. 4-4, 8). Exposures of the mation are unknown in the area.

The Sparta is about 300 feet thick in northern Madison County. An increase in ickness of about 300 feet occurs as the formation dips loward the Jackson dome from a northeast. The formation thins to less than 400 feet on the crest of the dome it thickens again southwest of the dome to more than 800 feet in southwestern Hinds anty. Thinning is apparent also over the Oakley salt dome and the Flora oil field, give 8.

#### Well Yields and Aquifer Characteristics

Yields of wells in the Cockfield Formation are as great as 500 gpm. Specific capacities of 10 gpm per foot have been measured. In general, specific capacities of wells in the Cockfield are less than those for wells in the Sparta.

Five aquifer tests were made in the Cockfield Formation (table 5). Transmissibility values ranged from 10,000 to 41,000 gpd per foot and the coefficient of permeability averaged 360 gpd per sq ft, considerably less than the average for the Sparta Sand. Permeability of Cockfield sand beds is highest east of Jackson. The effects of pumping from the Cockfield and a comparison of the Cockfield and Sparta aquifers are illustrated on the theoretical drawdown graph (fig. 11).

#### Water Quality

The Cockfield Formation contains hard water in the outcrop area of Madison County, but natural softening occurs down the dip. The iron content is high in the outcrop area, and even as far south as northern Hinds County it is as much as 3.5 pm. Total mineralization increases down the dip as the water changes from a calcium bicarbonate to a sodium bicarbonate type. The dissolved solids content ranges from about 200 ppm in Madison County to 1,310 ppm at Utica. Generally the fluoride content is low (table 4). However, water from well S3 at Utica contained 4.4 ppm of fluoride and vater from well D9, west of Edwards, contained 2.8 ppm. These are the only known occurrences, in the three counties, of water from the Cockfield having fluoride in excess of 1 ppm. The range in pH of the water was from 7.5 to 8.9, similar to that of water in the Sparta Sand.

North of the Jackson dome, little difficulty is experienced in obtaining clear water from the Jockfield. West and south of Jackson, coloring in the water is more of a problem and is indicative of more abundant lightle and longer contact of the water with lightlic materials.

#### Davelopment of Water Supplies

The Cockfield Formation is the source of more than helf of the municipal water supplies in the area (table 2). In the past, all supplies for the subdivisions around Jackson were developed in the Sparta Sand, but recently wells have been completed in the Cockfield west of the city. The Cockfield is unused in the city of Jackson: the term wells completed in the unit in past years having been abandoned because of fine sand and poor yield of the aquifer and the ready availability of water from the city system.

The heaviest concentration of Cockfield wells forms a belt 4 to 6 miles wide surrounding the Jackson dome. The Cockfield is potentially a useful aquifer under-lying the city of Jackson at shallow depths. Well yields and specific capacities protectly would be smaller than those of the Sparta, owing to the finer texture of the material composing the aquifer.

South and west of the main belt of wells in the Cockfield Formation the Forest Hill Sand and matahoula Sandstone are generally used for small-capacity wells, and only at places where neither of these units contains sufficient sand to yield a satisfactory where supply are wells drilled to the Cockfield Formation.

#### Alluvium

Alluvial deposits occur in small amounts in valleys of all the tributary streams and in greater volume in the valleys of the Big Black and Pearl Rivers. The deposits have a maximum thickness of about 40 feet in the central part of the valleys, of which more than one-half usually is saturated. The sources of water in the alluvium are precipitation in the valleys and underflow from adjacent uplands.

Deposits in the Pearl River valley consist of silt, sand, and pea gravel. They contain shell fragments, lignitic shale, and other materials derived locally by erosion of the Tertiary or younger beds. Coarser deposits prevail at the base of the unit and finer materials at the top. Where the Pearl River flows across the Yazoo Clay the valley is  $3\frac{1}{2}$  to 4 miles wide. Where the river crosses the Forest Hill, Vicksburg, and Catahoula cuestas on the south flank of the Jackson dome, the valley walls converge to a width of  $1\frac{3}{4}$  miles.

Domestic wells are common in the alluvium along the river valleys and furnish sufficient water for domestic and stock supplies. One small industrial water supply has been developed from the alluvium south of Jackson near the Pearl River.

Alluvium is limited in both valleys, but wells 6 inches in diameter probably are capable of producing 50 gpm or more. The permeability of the coarse part of the alluvial fill is estimated to be about 500 gpd per sq ft.

Water in the alluvium is a hard, calcium bicarbonate type containing 0.04 to 33 ppm of iron (table 4). Hardness as great as 341 ppm has been measured.

#### SUMMARY OF GROUND-WATER CONDITIONS

Water-bearing formations ranging in age from Eocene to Recent underlie the Jack son metropolitan area. The regional southwesterly dip of the formations is altered locally by the Jackson dome. Fresh water generally occurs to depths greater than 2,000 feet.

Five aquifers are available for development of moderate to large ground-water supplies. These are in the Wilcox Group, Sparta Sand, Cockfield Formation, Forest Hill Sand, and Catahoula Sandstone. Thick beds of sand in the Wilcox Group constitute a virtually untapped source of water supplies in the northeastern half of the area. The Sparta Sand is the most intensively developed of the aquifers. In 1960 an average of 11 mgd was withdrawn from the Sparta for industrial and municipal uses. Fresh water is available everywhere in the Cockfield Formation, and the unit supplies more than half of the small municipalities in the area. Most other water-supply developments in this aquifer consist of rural domestic supplies. The Forest Hill Sand is limited to the southern third of the project area. Thin beds of sand in this unit provide domestic supplies and a few small municipal supplies. Many domestic water supplies in the southern half of Hinds and Rankin Counties are obtained from the Catahoula Sandstone, as is the municipal supply at Utica.

Ground water in shallow zones is low in dissolved-solids content, generally is soft, and has a pH less than 7. Iron is a common objectionable feature in the water but it may be a product of corrosion of well fittings. In deeper zones the water is a soft, sodium bicarbonate type having a pH great than 7. Dissolved-solids content averages about 400 ppm for the aquifers younger than the Wilcox Group; the average is somewhat higher for the Wilcox. Iron content rarely is a problem in the deep well High fluoride concentrations are common in water from the Wilcox Group and Forest High

### U.S. EPA REGION IV

# **SDMS**

Unscannable Material Target Sheet

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	Castocing Cargoring							
Nature of Material:								
Map:	Computer Disks:							
Photos:	CD-ROM:							
Blueprints:	Oversized Report:							
Slides:	Log Book:							
Other (describe): <u>Healogic Map of Mississippi</u>								
Amount of material:								
* Please contact the appropriate Records Center to view the material *								

Reference 14
Gulf States Creosoting Company
Flowood; Rankin Co., Mississippi
MSN000407423

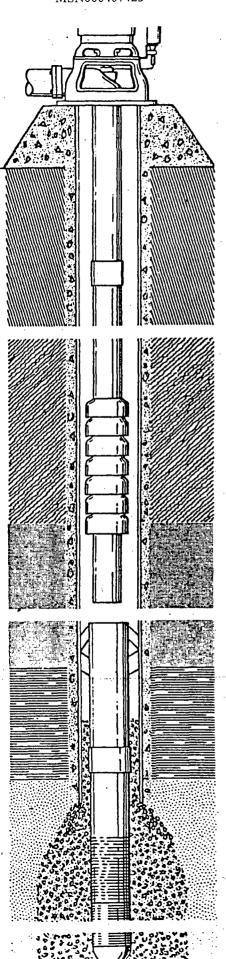
### GRINER DRILLING SERVISE

TELEPHONE 601/736-6347
POST OFFICE DRAWER 825
COLUMBIA, MISS. 39429

### LOG FORM

NAME Town of Flowood
LOCATION NE/4, NW/4, SE/4, NE/4, Section 12.
ENGINEER Lester Engineering Co.
DRILLER Allen Sistrunk LOG. NO K68
COMPLETED April 30 19 82 ACCEPTED
Sales Engineer T.N. SHOWS Field Supervisor F. SISTRUNK

COMPLETED April 30 19 82 ACCEPTED
Sales Engineer T.N. SHOWS Field Supervisor F. SISTRUNK
WELL DATA
Length surface casing N/A; size surface casing N/A
Cemented N/A No. Sacks N/A Size Drilled Hole 21"
Depth drilled hole 562 Size well casing 16" Type Welded
Length well casing 472' cemented Yes No. Sacks 600
Size underreamed hole 32" length underreamed hole 82'
Size screen 8" type Bar Weld mtg. by Houston Well Sc
Slot size .016 material 304 S.S. length screen 82'
Lap pipe size 8" lap pipe length 62' type Welded
Type gravel Ratcliff No. yds 12 Distance to lap 422
Distance to screen top 482 distance to gravel
Distance to screen top 562 time better Back Wash Valve
Distance to screen bottom 562 type bottom Back Wash Valve  Connection top of lap8" R&L Collar static water level 185.78
Connection top of rape
PUMP DATA  Type Turbine make Floway Serial No. 8220415
Size bowls 10" No. stages 12 Curve No DKH Length bowls 9'6"
Length column 280' size column 8" type column T & C
Size oil tube N/A size shaft 11 length suction 201
Size suction 811 size discharge 811 Head No C-16111x811x
Overall pump length 289'6" Length headshaft 10'3"
Type lubrication Water type oiler N/A length air line N/A
Rated capacity 750 GPM Total Head 385
RPM 1770 Size foundation 2' x 2' Height 2'
ELECTRIC MOTOR DATA
Type VHS Make G.E. Serial NOTJ422103
HP 125 Voltage 480 RPM 1800 Frame B405TP16
Style P Phase 3 Cycle 60 AMPS 144
Height motor 36½ Dia. base 16½ Clutch bore 1½ Clutch No
Top bearing No GOO1 Lower bearing No 93PO16 Lubrication Oil & Gre
SWITCH DATA
Type Pumping Plant Panel Make Control Systems
Catalog No N/A Size 4 Rating 125 H.P.
Volts 480 Phase 3 Cycles 60
HP rating 125 Size relays 3 Relay type Quick Trip
Entrance switch 200 Amp
Fuse size & type 225 Amp Breaker Size wire 3-0



DICK DERVICE, MINU

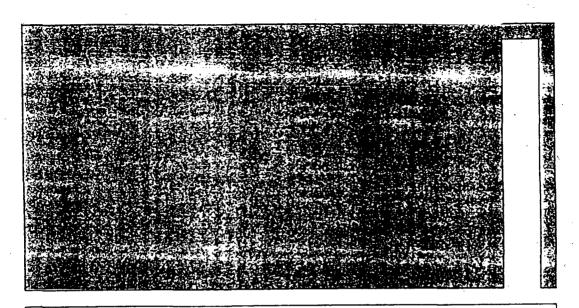
## TELEPHONE 601/736-6347 POST OFFICE DRAWER 825 COLUMBIA, MISS. 39429

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TEST HO		E LOCATION NEW NAME SET NEW Sec. 12	
TOTAL DEPTH	THICKNESS EACH STRATUM	FORMATION	REMARKS
3	3	Top soil	
39	36	Clay with thin sand streaks	
172	133	Clay	
213	41	Sand, fine	
444	231	Clay & rocks	
460	16	Sand & clay streaks	
468	8	Clay	•
570	102	Sand, medium	
604	34	Clay	
628	24	Sand	
661	33	Clay & sand streaks	
725	64	Sand with Thin clay streaks	
806	81	Clay & rocks	
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Date Starte	d 10-12-	19_81 Date Completed	10-16 19 81
. •		T. N. Shows	Field Engineer

Allen Sistrunk

\_\_Driller

Reference 15 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423



## R. Allan Freeze

Department of Geological Sciences University of British Columbia Vancouver, British Columbia

## John A. Cherry

Department of Earth Sciences University of Waterloo Waterloo, Ontario

## GROUNDWATER

Prentice-Hall, Inc. Englewood Cliffs, New Jersey 07632

Table 2.2 Range of Values of Hydraulic Conductivity and Permeability

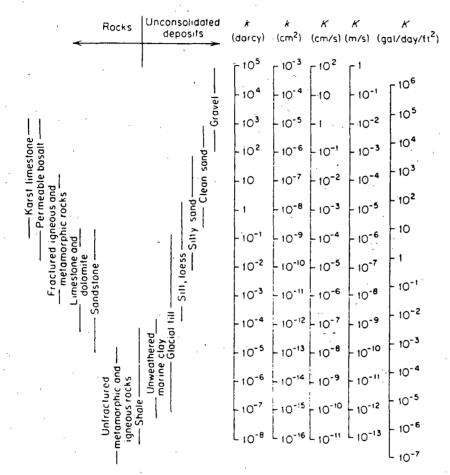


Table 2.3 Conversion Factors for Permeability and Hydraulic Conductivity Units

		Permeability, k*	<u> </u>	Hy	draulic conducti	vity, K
_	cm²	ft²	darcy	m/s	ft/s	U.S. gat/day/ft²
cm²	. 1	1.08 10-3	1.01 . 10*	9.80 - 102	3.22 × 10 <sup>3</sup>	1.85 < 109
ft <sup>2</sup>	9.29 / 102	1	$9.42 \cdot 10^{10}$	9.11 < 105	$2.99 \times 10^{6}$	$1.71 \times 10^{12}$
darcy	9.87 × 10-9	$1.06 \times 10^{-11}$	1	9.66 × 10-6	3,17 × 10~5	$1.82 \times 10^{1}$
m s	1.02 10-3	1.10 10-6	1.04 - 105	1	3.28	2.12 106
ft/s	3.11 × 10-4	3.35 × 10-7	$3.15 \times 10^{4}$	3.05 10-1	1	$6.46 \times 10^{5}$
U.S. gal da	ay ft <sup>2</sup> 5,42 × 10 <sup>-10</sup>	5.83 × 10-13	5.49 10	4.72 - 10-1	1.55 - 10-6	1

<sup>•</sup>To obtain k in ft<sup>2</sup>, multiply k in cm<sup>2</sup> by 1.08  $\times$  10<sup>-3</sup>.

Reference 16 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423

## Alexicological Profile For

## **CREOSOTE**

## U.S. DEPARTMENT OF HEALTH & HUMAN SERVICES

Public Health Service

Agency for Toxic Substances and Disease Registry

TRAJIEJO

## TOXICOLOGICAL PROFILE FOR CREOSOTE

#### Prepared by:

Clement International Corporation Under Contract No. 205-88-0608

Prepared for:

Agency for Toxic Substances and Disease Registry U.S. Public Health Service

December 1990

#### DISCLAIMER

The use of company or product name(s) is for identification only and does not imply endorsement by the Agency for Toxic Substances and Disease Registry.

#### FOREWORD

The Superfund Amendments and Reauthorization Act (SARA) of 1986 (Public Law 99-499) extended and amended the Comprehensive Environmental Response, Compensation, and Liability Act of 1980 (CERCLA or Superfund). This public law directed the Agency for Toxic Substances and Disease Registry (ATSDR) to prepare toxicological profiles for hazardous substances which are most commonly found at facilities on the CERCLA National Priorities List and which pose the most significant potential threat to human health, as determined by ATSDR and the Environmental Protection Agency (EPA). The lists of the 250 most significant hazardous substances were published in the Federal Register on April 17, 1987, on October 20, 1988, on October 26, 1989, and on October 17, 1990.

Section 104(i)(3) of CERCIA, as amended, directs the Administrator of ATSDR to prepare a toxicological profile for each substance on the list. Each profile must include the following content:

- (A) An examination, summary, and interpretation of available toxicological information and epidemiological evaluations on the hazardous substance in order to ascertain the levels of significant human exposure for the substance and the associated acute, subacute, and chronic health effects,
- (B) A determination of whether adequate information on the health effects of each substance is available or in the process of development to determine levels of exposure which present a significant risk to human health of acute, subacute, and chronic health effects, and
- (C) Where appropriate, an identification of toxicological testing needed to identify the types or levels of exposure that may present significant risk of adverse health effects in humans.

This toxicological profile is prepared in accordance with guidelines developed by ATSDR and EPA. The original guidelines were published in the <u>Federal Register</u> on April 17, 1987. Each profile will be revised and republished as necessary, but no less often than every three years, as required by CERCLA, as amended.

The ATSDR toxicological profile is intended to characterize succinctly the toxicological and adverse health effects information for the hazardous substance being described. Each profile identifies and reviews the key literature (that has been peer-reviewed) that describes a hazardous substance's toxicological properties. Other pertinent literature is also presented but described in less detail than the key studies. The profile is not intended to be an exhaustive document; however, more comprehensive sources of specialty information are referenced.

THE BUILDING TO YOUR

#### Foreword

Each toxicological profile begins with a public health statement, which describes in nontechnical language a substance's relevant toxicological properties. Following the public health statement is information concerning significant health effects associated with exposure to the substance. The adequacy of information to determine a substance's health effects is described. Data needs that are of significance to protection of public health will be identified by ATSDR, the National Toxicology Program (NTP) of the Public Health Service, and EPA. The focus of the profiles is on health and toxicological information; therefore, we have included this information in the beginning of the document.

The principal audiences for the toxicological profiles are health professionals at the federal, state, and local levels, interested private sector organizations and groups, and members of the public.

This profile reflects our assessment of all relevant toxicological testing and information that has been peer reviewed. It has been reviewed by scientists from ATSDR, the Centers for Disease Control, the NTP, and other federal agencies. It has also been reviewed by a panel of nongovernment peer reviewers and is being made available for public review. Final responsibility for the contents and views expressed in this toxicological profile resides with ATSDR.

William L. Roper, M.D., M.P.H.

Administrator

Agency for Toxic Substances and Disease Registry

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This Statement was prepared to give you information about creosote and to emphasize the human health effects that may result from exposure to it. The Environmental Protection Agency (EPA) has identified 1,177 sites on its National Priorities List (NPL). Creosote as been found at 31 of these sites. However, we do not know how many of the 1,177 NPL sites have been evaluated for creosote. As EPA evaluates more sites, the number of sites at which creosote is found may change. The information is important for you because creosote may cause harmful health effects and because these sites are potential or actual sources of human exposure to creosote.

When a chemical is released from a large area, such as an industrial plant, or from a container, such as a drum or bottle, it enters the environment as a chemical emission. This emission, which is also called a release, does not always lead to exposure. You can be exposed to a chemical only when you come into contact with the chemical. You may be exposed to it in the environment by breathing, eating, or drinking substances containing the chemical or from skin contact with it.

If you are exposed to a hazardous substance such as creosote, several factors will determine whether harmful health effects will occur and what the type and severity of those health effects will be. These factors include the dose (how much), the duration (how long), the route or pathway by which you are exposed (breathing, eating, drinking, or skin contact), the other chemicals to which you are exposed, and your individual characteristics such as age, sex, nutritional status, family traits, life style, and state of health.

#### 1.1 WHAT IS CREOSOTE?

Creosote is a complex mixture of many chemicals. There are three kinds of creosote. One type results from high-temperature treatment of coal (coaltar creosote), one results from high-temperature treatment of beech and other woods (beechwood ereesote), and one comes from the resin of the creosote bush (creosote bush resin). Coal-tar creosote is the most widely used wood preservative in the United States. Coal-tar products are also ingredients in medicines which are used to treat skin diseases. About 300 chemicals have been identified in coal-tar creosote, and there may be 10,000 other chemicals present in the mixture. The major chemicals in coal-tar creosote that can cause harmful health effects are polycyclic aromatic hydrocarbons (PAHs), phenol, and cresols. Beechwood creosote was at one time used as a disinfectant and as a treatment for coughs, but it is rarely used today. The major chemicals present in beechwood creosote are phenol, cresols, and guaiacol. Because coal-tar creosote is the only type found in the environment and hazardous waste sites, its effects on human health will be emphasized in

this toxicological profile. Furthermore, because this profile specifically addresses the creosote mixture, effects on health from exposure to the individual major chemicals in creosote, namely, the PAHs or phenol, will not be discussed in any great detail. The Agency for Toxic Substances and Disease Registry (ATSDR) Toxicological Profile for Polycyclic Aromatic Hydrocarbons and the ATSDR Toxicological Profile for Phenol provide more information on these chemicals.

Coal-tar creosote is usually a heavy, oily, liquid that is typically amber to brown in color. Mixtures of creosote and other coal-tar products are black. The creosote found at hazardous waste sites is most often a black, heavy liquid. It has a sharp smoky odor, and a burning taste. It burns easily, but does not dissolve readily in water.

Creosote does not occur naturally in the environment but it can be released to water and soil through its use as a wood preservative. However, there are many known natural sources for components of the creosote mixture, such as PAHs. For more information on the natural sources of these components, the reader can refer to the ATSDR Toxicological Profile for Polycyclic Aromatic Hydrocarbons. Some parts of the creosote mixture can enter groundwater or change into other substances while other parts persist in treated wood products for decades. More information on the chemical and physical properties of creosote can be found in Chapter 3 and on its occurrence and fate in the environment in Chapters 4 and 5.

#### 1.2 HOW MIGHT I BE EXPOSED TO CREOSOTE?

The major sources of human exposure to coal-tar creosote are contaminated hazardous waste sites, wood treatment facilities, and wood products treated with creosote. You cannot buy coal-tar creosote for treating wood products in your home. Wood products that are typically treated with creosote are railroad ties used by the railroads and for landscaping, telephone poles, marine pilings, and fence posts. You can also be exposed to creosote through contact with soil, water, or air contaminated as the result of releases from waste disposal sites and wood treatment facilities and the burning of treated scrap wood. Exposure of the general population to creosote from currently operating wood treatment facilities should be minimal because all of these facilities are subject to strict controls by the Occupational Safety and Health Administration (OSHA) and the Environmental Protection Agency (EPA). However, exposure to creosote from abandoned wood treatment facilities is possible. There should be very little exposure of the average homeowner to creosote solutions used for wood treatment because it can only be sold to certified applicators, however, you can still be exposed to creosote treated products.

There is no information on background levels of creosote in air or food. Background levels of certain parts of the creosote mixture such as PAHs can be found in some city soils and foods, but it is not known if these substances

came from creosote. For more information on human exposure to creosote see Chapter 5.

#### 1.3 HOW CAN CREOSOTE ENTER AND LEAVE MY BODY?

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od. be Creosote can enter your body through the lungs as a contaminant of air, through the stomach and intestines after eating contaminated food or drinking contaminated water, or through the skin. Although there is no information on how fast or how much of the creosote mixture is absorbed, many of the parts of the creosote mixture (for example, PAHs) are rapidly absorbed through the lungs and the stomach and the intestines.

The most common routes of exposure around hazardous waste sites are likely to be through the skin, and drinking water contaminated with creosote. Although it is not known how rapidly creosote can enter the body through the skin, creosote can cause reddening just from skin contact. Eating soil contaminated with coal-tar creosote can also provide a source of exposure. Chemicals in coal-tar creosote appear to accumulate in the body, particularly in fat tissue. Most of the chemicals in creosote that are taken into the body and are not stored in the body tissues, leave in the feces within a few days. More information on how creosote enters and leaves the body can be found in Chapter 2.

#### 1.4 HOW CAN CREOSOTE AFFECT MY HEALTH?

Reports describing coal-tar creosote poisoning in workers or accidental or intentional ingestion of coal-tar creosote indicate that brief exposures to large amounts of coal-tar creosote can cause harmful effects on your skin, eyes, nervous system, and kidneys and can result in death. Longer-term exposure to lower levels of coal-tar creosote can also result in damage to your skin, such as reddening, blistering or peeling.

The major organs or systems affected by longer-term exposure to lower levels of coal-tar creosote in animals are the skin and lungs, whereas only the skin has been observed to be affected in humans under these exposure conditions. All of these effects worsen as the level of coal-tar creosote exposure increases.

An increased risk for cancer has been demonstrated in animals exposed to coal-tar creosote. Birth defects have been seen in livestock exposed to coal-tar creosote-treated wood. Since these effects were seen in animals, it is also possible that they could occur in humans. However, we have no information about these effects in humans. More information on the health effects associated with exposure to creosote is presented in Chapter 2.

#### 1.5 WHAT LEVELS OF EXPOSURE HAVE RESULTED IN HARMFUL HEALTH EFFECTS?

Tables 1-1 through 1-4 show that no information is available on specific amounts of creosote in air, water, or food that cause health effects other than death. Coal-tar creosote can cause death to humans and animals if large enough quantities are swallowed. No information is available on the health effects of inhaled creosote.

Skin contact with a few drops of coal-tar creosote irritates and burns the skin and eyes. Coal-tar creosote also makes the skin more sensitive to the effects of the sun. These effects include burning, irritation and swelling.

### 1.6 IS THERE A MEDICAL TEST TO DETERMINE WHETHER I HAVE BEEN EXPOSED TO CREOSOTE?

There is no medical test to determine if you have been exposed to the creosote mixture. However, in your body, chemicals contained in creosote (PAHs) can attach to substances within the body. The presence of PAHs attached to these substances can be measured in body tissues or blood following exposure to creosote. However, this test is still being developed, and it is not known yet how well it works. In addition, chemicals contained in coal-tar creosote and their breakdown products can be measured in the urine of exposed individuals. Tests on the urine are commonly done to determine if a person has been exposed to coal-tar creosote. This test is not routinely available at a doctor's office and would require special equipment for sampling and detection of the compound. Although these tests can confirm that a person has been exposed to chemicals contained in coal-tar creosote, it is not yet possible to use the test results to accurately predict the severity of any health effects that might occur. Furthermore, these tests are not specific for the chemicals contained in coal-tar creosote, and the chemicals measured could have come from exposure to other sources. Additional information on tests for the measurement of creosote in the body is presented in Chapters 2 and 6.

## 1.7 WHAT RECOMMENDATIONS HAS THE FEDERAL GOVERNMENT MADE TO PROTECT HUMAN HEALTH?

The federal government has not developed regulatory standards and guidelines to protect people from the potential health effects of exposure to coal-tar creosote in drinking water and food. However, regulatory standards and guidelines in air and water exist for the key individual PAHs and phenols contained in creosote. For more information on the regulations and guidelines for PAHs and phenols, please refer to the ATSDR Toxicological Profiles for Polycyclic Aromatic Hydrocarbons and Phenol. EPA has declared creosote a restricted use pesticide which means that it can only be bought and used by certified applicators and only for those uses covered by the applicator's certification. The EPA has concluded that any release of creosote to the environment in excess of 1 pound should be reported.

TABLE 1-1. Human Health Effects from Breathing Coal-tar Creosote

Short-term Exposure (less than or equal to 14 days)

Levels in air (ppm)

Length of Exposure

Description of Effects
The health effects resulting
from short-term exposure of
humans to air containing
specific levels of creosote
are not known.

Long-term Exposure (greater than 14 days)

Levels in air (ppm)

Length of Exposure

The health effects resulting from long-term exposure of humans to air containing specific levels of creosote are not known.

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TABLE 1-2. Animal Health Effects from Breathing Coal-tar Creosote

Short-term Exposure (less than or equal to 14 days)

Levels in air (ppm)

Length of Exposure

Description of Effects
The health effects resulting
from short-term exposure of
animals to air containing
specific levels of creosote
are not known.

Long-term Exposure (greater than 14 days)

Levels in air (ppm)

Length of Exposure

Exposure of Effects

The health effects resulting from long-term exposure of animals to air containing specific levels of creosote are not known.

TABLE 1-3. Human Health Effects from Eating or Drinking Coal-tar Creosote

	Short-term Expo (less than or equal t	
Levels in Food (ppm)	Length of Exposure	Description of Effects The health effects resulting from short-term exposure of humans to food containing specific levels of creosote are not known.
<u>Levels in Water (ppm)</u>		The health effects resulting from short-term exposure of humans to water containing specific levels of creosote are not known.
	Long-term Expo (greater than 14	
Levels in Food (ppm)  Levels in Water (ppm)	Length of Exposure	Description of Effects The health effects resulting from long-term exposure of humans to food containing specific levels of creosote are not known.
ACTOR IN WALET (DDM)		The health effects resulting from long-term exposure of humans to water containing specific levels of creosote are not known.

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TABLE 1-4. Animal Health Effects from Eating or Drinking Coal-tar Creosote

Short-term Exposure (less than or equal to 14 days)

Levels in Food (ppm) Length of Exposure

Description of Effects

34,000

<24 hours

Causes death in rats

Levels sater (ppm)

The health effects resulting from short-term exposure of animals to water containing specific levels of creosote are not known.

Long-term Exposure (greater than 14 days)

els in Food (ppm) Length of Exposure

Description of Effects
The health effects resulting
from long-term exposure of

animals to food containing specific levels of creosote are not known.

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Levels in Water (ppm)

The health effects resulting from long-term exposure of animals to water containing specific levels of creosot are not known.

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The federal government has developed regulatory standards and guidelines protect workers from the potential health effects of other coal-tar dducts in air. OSHA has set a legally enforceable limit (threshold limit lue, or TLV) of 0.2 mg/m³ coal-tar pitch volatiles in workroom air to otect workers during an 8-hour shift.

For more information on regulations and advisories for coal-tar creosote posure, see Chapter 7.

#### 8 WHERE CAN I GET MORE INFORMATION?

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1g :e If you have any more questions or concerns not covered here, please contact your State Health or Environmental Department or:

Agency for Toxic Substances and Disease Registry Division of Toxicology 1600 Clifton Road, E-29 Atlanta, Georgia 30333

This agency can also give you information on the location of the nearest occupational and environmental health clinics. Such clinics specialize in recognizing, evaluating, and treating illnesses that result from exposure to hazardous substances.

#### HEALTH EFFECTS

#### INTRODUCTION

This chapter contains descriptions and evaluations of studies and interpretation of data on the health effects associated with exposure to creosote. Its purpose is to present levels of significant exposure for creosote based on toxicological studies, epidemiological investigations, and environmental exposure data. This information is presented to provide public health officials, physicians, toxicologists, and other interested individuals and groups with (1) an overall perspective of the toxicology of creosote and (2) a depiction of significant exposure levels associated with various adverse health effects.

#### 2.2 DISCUSSION OF HEALTH EFFECTS BY ROUTE OF EXPOSURE

To help public health professionals address the needs of persons living or working near hazardous waste sites, the data in this section are organized first by route of exposure -- inhalation, oral, and dermal -- and then by health effect -- death, systemic, immunological, neurological, developmental, reproductive, genotoxic, and carcinogenic effects. These data are discussed in terms of three exposure periods -- acute, intermediate, and chronic.

Levels of significant exposure for each exposure route and duration (for which data exist) are presented in tables and illustrated in figures. The points in the figures showing no-observed-adverse-effect levels (NOAELs) or lowest-observed-adverse-effect levels (LOAELs) reflect the actual doses (levels of exposure) used in the studies. LOAELs have been classified into "less serious" or "serious" effects. These distinctions are intended to help the users of the document identify the levels of exposure at which adverse health effects start to appear, determine whether or not the intensity of the effects varies with dose and/or duration, and place into perspective the possible significance of these effects to human health.

The significance of the exposure levels shown on the tables and graphs may differ depending on the user's perspective. For example, physicians concerned with the interpretation of clinical findings in exposed persons or with the identification of persons with the potential to develop such disease may be interested in levels of exposure associated with "serious" effects. Public health officials and project managers concerned with response actions at Superfund sites may want information on levels of exposure associated with more subtle effects in humans or animals (LOAEL) or exposure levels below which no adverse effects (NOAEL) have been observed. Estimates of levels posing minimal risk to humans (minimal risk levels, MRLs) are of interest to health professionals and citizens alike.

#### 2. HEALTH EFFECTS

For certain chemicals, levels of exposure associated with carcinogenic effects may be indicated in the figures. These levels reflect the actual doses associated with the tumor incidences reported in the studies cited.

Estimates of exposure posing minimal risk to humans (MRLs) are made, where data are believed reliable, for the most sensitive noncancer end point for each exposure duration. MRLs include adjustments to reflect human variability and, where appropriate, the uncertainty of extrapolating from laboratory animal data to humans. Although methods have been established to derive these levels (Barnes et al. 1987; EPA 1986a), uncertainties are associated with the techniques. The data on creosote are inadequate to calculate MRLs.

Creosote is a complex mixture of organic substances. It can be derived from three sources: coal-tar distillates (referred to herein as coal-tar creosote), wood (referred to herein as beechwood creosote), and the resin fr leaves of the creosote bush (<u>Larrea</u>, referred to herein as creosote bush resin). Coal-tar creosote is defined by the American Wood Preserver's Association (1988) as:

A distillate derived from coal-tar. As used in the wood preserving industry, creosote denotes a distillate of coal-tar produced by the high temperature carbonization of bituminous coal. Creosote consists principally of liquid and solid aromatic hydrocarbons and contains some tar acids and tar bases; it is heavier than water and has a continuous boiling range beginning at about 200°C.

Coal-tar creosote is widely used as a wood preservative, and is the for of creosote found in the environment and hazardous waste sites.

Beechwood creosote consists mainly of phenol, cresols, guaiacol, xylenol, and creosol. It is a colorless or pale yellowish liquid and it has characteristic smoky odor and burning taste (Miyazato et al. 1981). It had therapeutic applications in the past as a disinfectant and a stimulating expectorant, but it is rarely used today. Though beechwood creosote and coatar creosote have some components in common (e.g., phenols), and some of the adverse effects associated with exposure to beechwood creosote may be due to the phenol component, it is not known whether coal-tar creosote will induce these same effects. Furthermore, coal-tar creosote contains PAHs, some of which are carcinogenic, and beechwood creosote does not. Thus, the relevance of health effects data on beechwood creosote to risk associated with exposure to coal-tar creosote is not known.

Creosote bush resin consists of phenolics (e.g., flavonoids and nordihydroguaiaretic acid), neutrals (e.g., waxes), basics (e.g., alkaloids) and acidics (e.g., phenolic acids). The phenolic portion comprises 83%-91% the total resin. Nordihydroguaiaretic acid accounts for 5%-10% of the dry

#### 2. HEALTH EFFECTS

weight of the leaves (Leonforte 1986). Again, it is not known if the health effects associated with creosote bush resin are due to the phenolic components common to coal-tar creosote or if these effects would be expected to occur following exposure to coal-tar creosote.

The chemical composition of each of these forms of creosote can vary considerably, depending on the source of coal, wood, or plant, and the design and attendant operating conditions (temperature, gas distillation systems, etc.) used to produce the creosote. Throughout this profile, every attempt will be made to specify the characteristics of the creosote under discussion, and indicate which effects may be expected to be common to two or more forms. Furthermore, since the intent of this profile is to discuss the creosote mixture, and even though it is likely that creosote's toxicity is due largely to the major individual components, the PAHs and phenol, the health effects of these individual components will not be discussed in great detail. However, it is understood that the toxicity of the individual components may not be representative of the actual toxicity of the creosote mixture because of the possibility of synergistic and/or antagonistic interactions in the mixture. For more information on the health effects of these components, the reader can refer to the ATSDR Toxicological Profiles for Phenol and Polycyclic Aromatic Hydrocarbons (ATSDR 1989b, 1990).

#### 2.2.1 \ Inhalation Exposure

#### 2.2.1.1 Death

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No studies were located regarding death in humans following inhalation exposure to creosote. No deaths were observed in rats exposed to near saturated vapors of coal-tar creosote for one hour (concentration estimated to be less than 0.033 mL/L) (Pfitzer and Gross 1964). The exposed animals showed signs of slight eye and nose irritation and slight dyspnea. All rats were reported to exhibit weight gains comparable to the control animals and there were no treatment-related lesions observed at necropsy.

No studies were located regarding the following health effects in humans or animals following inhalation exposure to creosote.

- 2.2.1.2 Systemic Effects
- 2.2.1.3 Immunologic Effects
- 2.2.1.4 Neurologic Effects
- 2.2.1.5 Developmental Effects
- 2.2.1.6 Reproductive Effects

#### HEALTH EFFECTS

#### 2.2.1.7 Genotoxic Effects

No mutagenic activity was detected in the urine samples of workers briefly exposed to coal-tar creosote vapors in a wood preserving factory (Bos et al. 1984a). The frequency and level of exposure to creosote vapors was no reported. Of the three workers tested, two moved wood in and out of the wood impregnating cylinder and one operated the cylinder. All had the potential for brief inhalation exposure to creosote. Furthermore, the possibility of dermal exposure by contact with residual surface creosote cannot be excluded. The absence of mutagens in the urine samples tested by Salmonella typhimurium assay was attributed by the authors to a relatively low level of exposure, improper timing for urine sample collection, and the insensitivity of the assay. The data are not sufficient to draw a conclusion regarding the genotoxic potential of inhaled creosote vapors in humans.

No studies were located regarding genotoxic effects in animals followin, inhalation exposure to creosote.

#### 2.2.1.8 Cancer

Cases of multiple myeloma were compared in a Swedish case-referent study on past occupational and radiation exposure (Flodin et al. 1987). Exposure was assessed by means of a questionnaire mailed to the study subjects that asked questions concerning occupational exposures (including exposure to coal tar creosote) and radiation exposure. Radiation exposure included estimates. of natural background levels and estimates from medical procedures. exposure to gamma radiation, either through occupation, medical testing, or from background sources found in building materials, was not significantly associated with an increase in multiple myeloma. A significant crude rate ratio for occupational exposure to coal-tar creosote was obtained (6.0; p=0.001). The rate ratio point estimate for occupational exposure to creosote increased to 9.5 after stratifying for age. This result indicates an association between occupational creosote exposure and the subsequent development of multiple myeloma. This study was limited by the number of smokers in the case group as compared with the referent group, by gender differences between the cases and referents, and, possibly, by recall bias and incomplete case ascertainment. Furthermore, it is assumed that the exposure route was primarily dermal, although some inhalation exposure may have been possible. This study does suggest an area for further research, indicating a possible relationship between creosote exposure and the development of multiple myeloma.

Creosote does contain mutagenic PAHs, and therefore an association between creosote exposure and the development of multiple myeloma appears to be biologically feasible. Furthermore, creosote exposure has been associated with cancers of the nasal cavity, larynx, lung, skin, and scrotum (Fraumeni 1975). See Section 2.2.3.8, "Dermal Exposure, Cancer" for a more detailed,

#### 2. HEALTH EFFECTS

sion of the types of cancer associated with exposure to creosote in

No definitive inhalation carcinogenicity bioassays have been conducted received in animals.

#### Oral Exposure

Figure 2-1 and Table 2-1 describe the health effects observed in laboratory animals associated with oral exposure levels to coal-tar and beechwood creosote at varying time and exposure levels. Although beechwood creosote, creosote bush resin, and coal-tar creosote have some components in recommon (e.g., phenols), and some of the adverse effects associated with exposure to beechwood creosote may be due to the phenol component (e.g. acute and/or subacute toxicity), it is not known whether coal-tar creosote will induce these same effects. Furthermore, coal-tar creosote contains a complex mixture of carcinogenic/co-carcinogenic PAHs that probably account for the cancer risk associated with chronic exposure, and the other forms of creosote do not.

#### 2.2.2.1 Death

A 70-year-old man died following ingestion of an unspecified amount of "industrial" creosote (presumably coal-tar creosote) (Bowman et al. 1984). Death was attributed to multi-organ failure and occurred 30 hours after admission to the hospital. It is not known if this man had a history of prior creosote ingestion. Death has been reported to occur in adults and children 14 to 36 hours after the ingestion of about 7 g or 1-2 g coal-tar creosote, respectively (Clayton and Clayton 1981). The latter anecdotal information is of limited value because it was obtained from a secondary source that provided no supporting documentation. Thus, ingestion of creosote can be fatal to humans, but the dose level required to produce death cannot be accurately estimated from these reports.

Cases of lethal poisoning resulting from ingestion of large amounts of coal-tar creosote have been reported in larger farm animals. The acute fatal dose for sheep is 4-6 g/kg and for calves, over 4 g/kg (HSDB 1988). The oral LD<sub>50</sub> for coal-tar creosote is reported to be 433 mg/kg in the mouse and 725 mg/kg in the rat (RTECS 1981). However, another study reported an acute oral LD<sub>50</sub> of 1700 mg/kg in the rat (Pfitzer and Gross 1964). The reasons for this discrepancy in LD<sub>50</sub> values are not known. Based on these data, coal-tar creosote can be classified as mildly to moderately toxic.

The acute toxicity of beechwood crossote in both rats and mice was studied following the administration of single gavage doses in a 10% aqueous solution (Miyazato et al. 1981). The oral  $\rm LD_{50}$  of beechwood crossote in rats was 870 mg/kg (females) and 885 mg/kg (males). The highest dose at which no death occurred was 600 mg/kg. There was no significant difference between

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HEALTH EFFECTS

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	,		Exposure	: '		-	LOAE	L (Effec	t)		
Figure Key	Species	Route	Frequency/ Duration	Effect	NOAEL (mg/kg/da	у)	Less Serious (mg/kg/day)		Serious (mg/kg/day)	Reference	Form
ACUTE EXI	POSURE		<del></del>	:							
Death				. !	`						
1	Ret	(G)	l d lx/d					1700	(LD50)	Pfitzer and Gross 1964	coal
2	Rat	(G)	1 d lx/d	. !	600				(LD50). (LD50)	Miyazato et al. 1981	beech
3 .	Mouse	(G)	1 d 1x/d		M:376 F:313				(LD50) (LD50)	Miyazato et al. 1981	beech
Systemic	;			:							
4	Rat	(G)	1 d 1x/d	Gastro	•	2.52	(hyperemia, distension			Pfitzer and Gross 1964	coal
							of stomach)		•	٠.	
Reurolog	gical			· .							
5	Rat	(G)	1 d 1x/d				•	600	(convulsions)	Miyazato et al. 1981	beech
6	Mouse	(G)	1 d 1x/d					313	(convulsions)	Miyazato et al. 1981	beech
INTERMEDI	IATE EXPOSI	JRE		1					٠.		
Death				:			· .				
7 ·	Rat	(F)	3 mo 7d/wk		M: 1224	·			e.	Miyazato et al. 1981	. beech
			•		F: 768						
8	Mouse	(F)	3 mo 7d/wk		M: 1810					Miyazato et al. 1981	beech
				:	F: 1570					,	
Systemic	:	•	٠				•				
9	Rat	(F)	3 mo 7d/wk	Hemato			(decreased red blood cell, in-			Miyazato et al. 1981	beech
	./			· !			creased serum cholesterol)	•			

TABLE 2-1. Levels of Significant Exposure to Creosote - Oral

# TABLE 2-1 (Continued)

				Exposure				, LOAEL (E	ffect)	•	j r
Figure Key	Species		Frequency/ Duration	Effect	NOAEL (mg/kg/da		Less Serious (mg/kg/day)	Serious (mg/kg/day)	Reference	Form	
•			·	Hepatic			(increased liver weight:body			· .	
		. •		Renal	-		weight ratio) (increased kidney weight:body				
	-			Other		1224	ratio) (decreased body weight and food intake)	•			
				Other		M:	(increased spleen body weight ratio)				
10	Mouse	(F)	3 mo 7d/wk	Hemato	M: 1810 F: 1570				Miyazato et al. 1981	beech	
		-		Hepatic	. 1370	F:345	(increased liver weight:body weight ratio)	•			
			1	Renal			(decreased kidney weight)	;			
				Other		1810	(decreased body weight and food intake)				
Neurolog	gical				•			•	•		
11 • .	Rat	(F)	3 mo 7d/wk			M: 257	(increased brain weight;body weight ratio)	•	Miyazato et al. 1981	beach	
Reproduc	ctive										
12	Rat	(F)	3 mo 7d/wk		•		(increased testis: body weight ratio)		Miyazato et al. 1981	beech	
CHRONIC E	EXPOSURE										
Death											
13	Rat	(F)	96 wk 7d/wk		•	٠.		M:313 (death)	Miyazato et al. 1984b	beech	

		·							· · · · · · · · · · · · · · · · · · ·	
•			Exposure				LOAEL (Ef:	(ect)		
Figure Key	Species		Frequency/ Duration	Effect	NOAEL (mg/kg/day		Less Serious (mg/kg/day)	Serious (mg/kg/day)	 Reference	Form
Systemic				<del>-                                    </del>			<del></del>			<u> </u>
14	Rat	(F)	96 wk 7d/wk	Hemato Hepatic	1	F:179	(increased serum cholesterol) (increased liver weight:body weight ratio)		Miyazato et al. 1984b	beech
				Renal			(increased kidney weight:body weight ratio; in- creased BUN (male)			
	:			Other	M:313 F:394		nephrosis (mele) (decresed food intake)			
		٠		ļ. ļ			(decreased body weight)			
15	Mous •	(F)	52 wk 7d/wk	Resp Hemato		M: 247 P: 297	(histopath)		Miyazato et al. 1984a	beech
				Hepatic Renal	F:297		(increased liver weight:body weight ratio) (increased kidney			
	:			Other		1:247	weight:body weight ratio) (decreased body weight and food	• •		
Neurolog	ical			!			intake)		•	
16	Rat	(F)	96 wk 7d/wk	:	•	7:394	(increased brain weight:body weight ratio)		Miyazato et al. 1984b	beach
17	Mouse	(F)	52 wk 7d/wk		F		(increased brain weight:body weight ratio)		Miyazato et al. 1984a	beech

<sup>\*</sup>Converted to an equivalent concentration in food for presentation in Table 1-4.

mg/kg/day = milligrams per kilogram per day; d = day; mo = month; x/d = times per day; (G) = gavage; (F) = feed; NS = not specified; Gastro = gastrointestinal; Hemato = hematological; Resp = respiratory; BUN = blood urea nitrogen; LD50 = lethal dose, 50% kill;

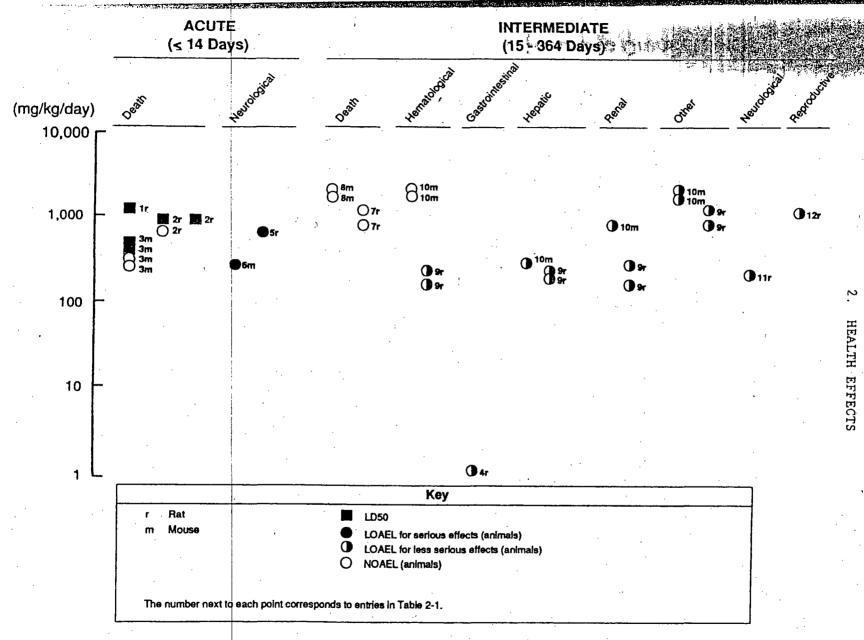


FIGURE 2-1. Levels of Significant Exposure to Creosote - Oral

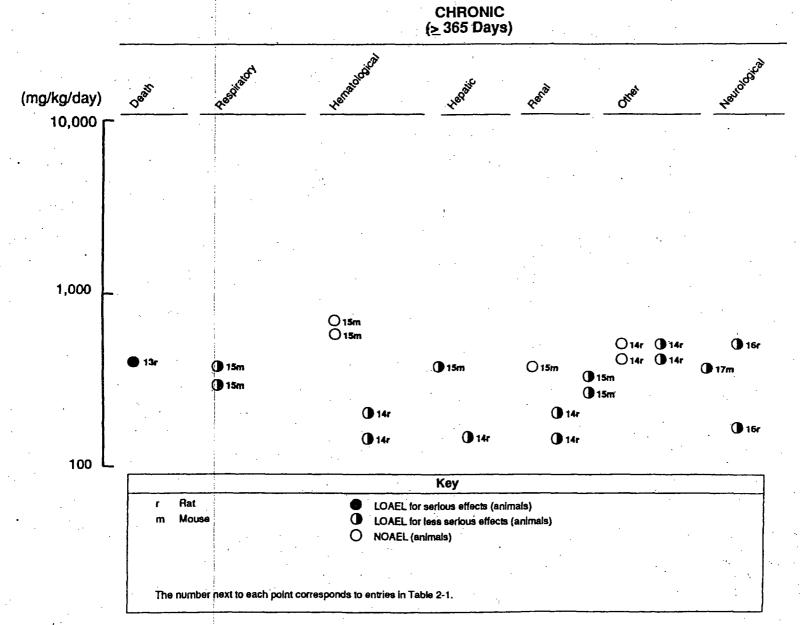


FIGURE 2-1. Levels of Significant Exposure to Creosote - Oral

(Continued)

mal and female rats with respect to mortality, and most animals died within 60 minutes following severe convulsions. Mice appeared to be more susceptible to the lethal effects of beechwood creosote. The oral LD50s in mice were 43 mg/kg (female) and 525 mg/kg (male). The highest dose at which no death coursed was 313 mg/kg (female) and 376 mg/kg (male). The mortality in female mice was significantly higher than in male mice. Most animals died within 30 minutes following severe convulsions. These results indicate that the acute toxicity of beechwood creosote is relatively low, and some species and sex differences exist.

The highest NOAEL values and all LD<sub>50</sub> values for death in each species are recorded in Table 2-1 and plotted in Figure 2-1.

# 2.2.2.2 Systemic Effects

There is relatively little information available regarding the systemic effects of ingested creosote. The database consists primarily of old anecdotal reports or animal studies that would be considered inadequate by current standards. Three recent studies published by Miyazato et al. (1981, 1984a,b) that evaluated the acute, subchronic, and chronic effects of beechwood creosote in rats and mice comprise the bulk of reliable information on the systemic effects of ingested creosote. Based on the results of these studies, the liver, kidney, and central nervous system may be target organs of creosote toxicity. Effects have also been observed in the gastrointestinal system, respiratory system, and cardiovascular system. No studies were located regarding musculoskeletal or dermal/ocular effects in humans or animals after oral exposure to creosote.

Respiratory Effects. No studies were located regarding respiratory effects in humans following oral exposure to creosote.

A slightly higher incidence of bronchitis or thickening of the tracheal mucous membrane was observed in mice who ingested feed that contained 0.3% (equivalent to 247 mg/kg/day for males and 297 mg/kg/day for females) and 0.6% (equivalent to 474 mg/kg/day for males and 532 mg/kg/day for females) beechwood creosote for 52 weeks (Miyazato et al. 1984a). The authors attributed this to irritation from long-term inhalation exposure to volatile components of creosote in the feed, and not a direct toxic effect on the respiratory tissue. These values are presented in Table 2-1 and plotted in Figure 2-1.

Beechwood creosote has been and continues to be used therapeutically on a limited basis as an expectorant/cough suppressant based on its presumed ability to increase the flow of respiratory fluids. The efficacy of creosote (type not specified, but presumably beechwood creosote) as an expectorant was studied by measuring the output of respiratory tract fluids in cats given a single dose of 0.1 or 5 mL/kg (concentration not specified) (Stevens et al.

1943). Creosote produced a slight increase in the output of respiratory trac fluid under these conditions. This is not considered a toxic effect. Given the limitations of this study (e.g., no dose information, no other respiratory effects evaluated), it provides no useful information on the potential respiratory effects of beechwood creosote.

Cardiovascular Effects. The case of a 52-year-old woman who had been taking creosote (type and dose not specified) for 9 years to treat chronic bronchitis was reported by Robinson (1938). The woman was found to be weak, dizzy, light-headed, and hypertensive (blood pressure = 206/140). A modified diet and diuretic therapy relieved all of these symptoms. Upon reinstitution of creosote therapy, her blood pressure rose to 235/130. The author concluded that creosote was responsible for the woman's hypertension. This study provides anecdotal evidence of creosote-induced cardiovascular effects, but the limited sample size, lack of detail on exposure, and possibility of confounding factors limits its usefulness.

No studies were located regarding the cardiovascular effects in animals following oral exposure to creosote.

Gastrointestinal Effects. Ulceration of the oropharynx and petechial hemorrhages over the gastrointestinal serosal surfaces were noted at autopsy in the case of a 70-year-old man who died following ingestion of industrial creosote (Bowman et al. 1984). However, the esophagus and stomach were intact. The authors attributed these effects to acute tissue damage resulting from phenol-induced corrosive effects (phenol is a component of creosote).

Animals that died following the administration of single gavage doses of coal-tar creosote in an acute range-finding study (doses ranged from 2.52-5.00 mg/kg) exhibited hyperemia and distention of the stomach upon necropsy. No studies were located regarding gastrointestinal effects in animals following intermediate- or chronic-duration oral exposure to creosote.

Hematologic Effects. No studies were located regarding hematologic effects in humans following oral exposure to creosote.

Various serum chemistry and hematological parameters were measured by Miyazato et al. (1981, 1984a,b) in rats and mice fed beechwood creosote in the daily diet for 3 months, 52 weeks (mice) or 96 weeks (rats). No significant treatment-related changes were noted in mice of either sex fed up to 1,570 (female) or 1,810 (male) mg/kg/day for 3 months (Miyazato et al. 1981). These doses are considerably higher than the oral LD<sub>50</sub>s reported for mice by the same authors. One possible explanation for this discrepancy is that the LD<sub>50</sub>s were determined by bolus gavage injections and in the subchronic study, the beechwood creosote was administered in the feed. A slight reduction in red blood cells (RBCs) and a slight increase in serum cholesterol was noted in male and female rats following dietary exposure to 163 (female) or 257 (male)

mg/rg/day and higher for 3 months (Miyazato et al. 1981). The reduction in RBOS was not dose-dependent and was therefore not considered to be foricologically significant; the significance of the cholesterol changes is not known.

Increases in serum cholesterol were also noted in rats exposed to creosote in the diet for 96 weeks at doses of 143 (male) and 179 (female) mg/kg/day and above (Miyazato et al. 1984b). Chronic (52 weeks) dietary exposure of mice to beechwood creosote resulted in statistically significant dose-related differences in mean cell volume, mean corpuscular hemoglobin, and absolute lymphocyte and neutrophil counts when compared to the corresponding control values. These changes were not considered by the authors to be toxicologically significant; they claimed that the values were within normal physiological ranges (Miyazato et al. 1984a).

The highest NOAEL values and all reliable LOAEL values for hematological effects in each species and duration category are recorded in Table 2-1 and plotted in Figure 2-1.

Hepatic Effects. Degeneration and necrosis of hepatocytes were observed at autopsy in the case of a 70-year-old man who ingested "industrial" creosote (Bowman et al. 1984). Given the advanced age of this man and the lack of comparison data, it is not possible to definitively attribute these effects to creosote ingestion.

Liver-to-body-weight ratios tended to increase in rats and mice exposed to beechwood creosote in the diet for either 3 months (rats and mice), 52 weeks (mice) or 96 weeks (rats) (Miyazato et al. 1981, 1984a,b). However, this response is of questionable toxicological significance, since no treatment-related changes were noted during histopathological evaluation. A significant increase in serum glutamic-oxaloacetic transferase (GOT) and glutamic-pyruvic transferase (GPT) levels was also observed in the chronically exposed female mice, but these levels were still within normal physiological range. Taken together, these early changes indicate that if increased doses of beechwood creosote are used, more clear cut hepatotoxic effects would be expected to occur. However, the liver weight changes observed may simply reflect pharmacological changes (e.g., response to enzyme induction).

All reliable LOAEL values for hepatic effects in each species and duration category are recorded in Table 2-1 and plotted in Figure 2-1.

Renal Effects. A 70-year-old man who ingested a fatal dose of "industrial" creosote became acidotic and anuric before he died, indicating the possibility of kidney failure (Bowman et al. 1984). Acute renal tubular necrosis was revealed at necropsy. However, the acute tubular necrosis may have been due to vascular insufficiency rather than a direct toxic effect on the kidney.

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Kidney-to-body-weight ratios tended to increase in rats and mice expose to beechwood creosote in the diet for either 3 months (rats and mice), 52 weeks (mice) or 96 weeks (rats) (Miyazato et al. 1981, 1984a,b). However, this response is of questionable toxicological significance in the subchronically treated mice and rats of both sexes and the chronically treate female rats since no treatment-related changes were noted at histopathologica evaluation. Blood urea nitrogen and serum inorganic phosphorus were elevated in the chronically treated male rats, which is indicative of uremia. Chronic nephrosis, which occurs spontaneously in old male rats, was also observed at higher incidence in the chronically treated male rats than in the control mal rats, and most likely accounts for the biochemical changes observed. The authors concluded that long-term exposure to beechwood creosote accelerated the occurrence of chronic nephrosis in male rats (Miyazato et al. 1984b). These results suggest that beechwood creosote has the potential to induce adverse effects in the kidney.

All reliable LOAEL values for renal effects in each species and duration category are recorded in Table 2-1 and plotted in Figure 2-1.

# 2.2.2.3 Immunological Effects

No studies were located regarding immunologic effects in humans or animals following oral exposure to creosote.

# 2.2.2.4 Neurological Effects

No studies were located regarding neurological effects in humans following oral exposure to creosote.

The first sign of poisoning following the gavage administration of single high doses (doses not specified) of beechwood creosote to mice and rats was muscle twitching followed by convulsions within 1 to 2 minutes. This was followed by asphyxiation, coma, and death (Miyazato et al. 1981).

Several cases of acute poisoning in cattle have been attributed to ingestion of coal-tar creosote. Six cattle believed to have licked creosote-treated electrical light poles as evidenced by burning over the mucosa of the mouth, tongue, and lips showed the following symptoms: extremely rapid respiration, contracted pupils, cold skin, apparent severe pain, and coma (Hanlon 1938). However, the possibility exists that some of these effects may have been due to ingestion of pentachlorophenol. Pentachlorophenol, like creosote, is an oil-borne wood preservative with extensive use in the public utility industry for treatment of utility poles. Some of the effects observed by Hanlon (1938) are more compatible with the metabolic effects (i.e., uncoupling of phosphorylative oxidation) associated with pentachlorophenol (for more information on the effects of pentachlorophenol, please refer to the ATSDR Toxicological Profile for Pentachlorophenol, ATSDR 1989a). Thus, it is

Epossible to determine conclusively whether creosote is toxic to the stral nervous system of animals based on this report.

Dose-related increased brain-to-body-weight ratios were observed in male ats exposed to beechwood creosote in the diet for 3 months (Miyazato et 1981), female mice exposed for 52 weeks (Miyazato et al. 1984a) and male and female rats exposed for 96 weeks (Miyazato et al. 1984b). This response is of questionable toxicological significance because of the lack of a dose-response trend and/or the lack of treatment-related pathological findings at necropsy.

All reliable LOAEL values for neurological effects for each species and duration category are recorded in Table 2-1 and plotted in Figure 2-1.

## 2.2.2.5 Developmental Effects

No studies were located regarding developmental effects in humans or animals following oral exposure to creosote.

# 2.2.2.6 Reproductive Effects

No studies were located regarding reproductive effects in humans following oral exposure to creosote. An increase in relative testis weight was observed in rats administered 1,224 mg/kg/day beechwood creosote in the diet for 3 months (Miyazato et al. 1981). There were no accompanying gross or histopathological lesions of the testes in these animals.

# 2.2.2.7 Genotoxic Effects

No studies were located regarding genotoxic effects in humans or animals following oral exposure to creosote.

### 2.2.2.8 Cancer

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No studies were located that dealt directly with an association between cancer and ingested creosote in humans. Excess cases of breast cancer have been observed in St. Louis Park, Minnesota that were tentatively associated with coal-tar creosote contamination of the water supply. Coal-tar creosote-derived PAHs were first detected in the water supply of St. Louis Park in November 1978, but may have been there for decades. A 100 acre plot of creosote-contaminated soil upon which stood a plant that used creosote and operated from 1917-1972 is believed to be the source of contamination. The levels of creosote or creosote-derived PAHs in the contaminated drinking water of St. Louis Park were not specified. There were 113 cases of breast cancer in St. Louis Park compared to 78 cases in the metropolitan Minneapolis-St. Paul area during 1969 to 1971. An attempt was made to demonstrate that these excess cases could be explained by known risk factors for breast cancer, such as age at first birth, parity, age at menarche and menopause, body mass index,

history of benign breast disease, and familial history and had no association with environmental exposure variables (Dean et al. 1988). The authors presented a method to adjust the breast cancer morbidity in St. Louis Park using data from a larger population with documented risk factors for breast cancer ("standard population"). These more stable rates based on breast cancer risk factors were determined in a larger study conducted by Helmrich et al. (1983). Dean et al. (1988) determined that the attributable risk due to the risk factors was higher for the breast cancer cases in St. Louis Park (0.598) than in the metropolitan Minneapolis-St. Paul area (0.311). They then used these attributable risks to calculate an adjusted morbidity ratio to estimate expected numbers of breast cancer cases in the community. After this adjustment, there appeared to be a higher expected number of breast cancer cases (n=134) than the observed number (113) in St. Louis Park, thereby negating any association with creosote in the water supply. It is necessary when using any standardization method in adjusting rates to ensure comparability among the groups in the study population that are being compared and between the study population and the standard population. It would appear that differences do exist between women in St. Louis Park, the metropolitan Minneapolis-St. Paul area, and the larger cohort studied by Helmrich et al. (1983). These differences were not thoroughly examined by Dean et al. (1988) and could include differences in demographic, economic, and/or environmental factors. However, Dean et al. (1988) did note a difference in religious backgrounds between the study population described by Helmrich (1983) and that of the St. Louis Park area. These dissimilarities indicate that the populations are not directly comparable and therefore the adjustment of rates is not appropriate. The authors did not attempt to incorporate creosote contamination of the water as an independent variable in their analysis.

The Minnesota Department of Health (1985) also reviewed the St. Louis Park data and concluded that this study did not provide adequate evidence to associate breast cancer with creosote-contaminated groundwater. They supported this conclusion with the following observations. It is not possible to classify individuals or residences within St. Louis Park according to their relative degree of historical exposure to PAH contaminants in drinking water because the pattern and history of municipal well contamination are not known; contaminant levels were measured at the well head and not at the tap; water treatment, storage, and distribution effects on contaminant concentration are not known; and much of the water distribution system is lined with coatings made of coal-tar or asphalt. Furthermore, given the ubiquitous nature of PAHs, it is probable that exposures to PAHs from food would significantly exceed exposures from contaminated St. Louis Park well water. In addition, it was found that the specific PAHs that have been shown to induce mammary tumors In rodents were either not present in contaminated wells or were detected very rarely even in the most highly contaminated wells, and the many published case-control and cohort studies of breast cancer have not demonstrated clearcut evidence of an association between breast cancer and smoking, which is a significant source of exposure to PAHs. These studies have also identified a number of risk factors that account for some of the observed variations in

Tites among different groups of women, and the women of St. Louis Park differ those in the general Metro area with respect to several of the factors had are known to influence breast cancer rates.

Dietary exposure of male and female mice to beechwood creosote at concentrations up to 532 mg/kg/day (which induced signs of toxicity) for \$2 weeks induced no treatment-related increase in the incidence of tumors (Miyazato et al. 1984a). This study is limited in that 52 weeks may not be a sufficient treatment duration to observe an increase in the incidence of fumors in mice. However, these same authors reported that dietary exposure of male and female rats to 394 mg/kg/day beechwood creosote (which induced signs of toxicity) for 96 weeks also failed to result in any treatment-related increase in the incidence of tumors (Miyazato et al. 1984b). Based on these lesults, there is no evidence that ingested beechwood creosote is carcinogenic to mice or rats.

## 2.2.3 Dermal Exposure

Table 2-2 describes the health effects observed in laboratory animals associated with dermal exposure to coal-tar creosote. Coal-tar creosote exerts its toxic effects primarily via dermal exposure. Several reports were found which describe the occurrence of dermal and ocular irritation, burns, and "warts" (i.e. squamous papillomas) following acute or prolonged skin contact with coal-tar creosote. Coal-tar creosote also induces phototoxicity of the skin, and has been demonstrated to be a skin carcinogen in animals.

### 2.2.3.1 Death

No studies were located regarding death in humans following dermal exposure to creosote. The dermal  $LD_{50}$  in rabbits has been estimated to be greater than 7.95 g/kg following a 24-hour application of coal-tar creosote to both intact and abraded skin (Pfitzer and Gross 1964).

# 2.2.3.2 Systemic Effects

The skin is the target organ for creosote toxicity following dermal exposure. No studies were located regarding respiratory, cardiovascular, hematologic, musculoskeletal, hepatic, or renal effects in humans or animals following dermal exposure to creosote.

Gastrointestinal Effects. No studies were located regarding gastrointestinal effects in humans following dermal exposure to creosote. Rabbits that died following single dermal applications of undiluted coal-tar creosote exhibited hyperemia of the intestines (Pfitzer and Gross 1964).

Dermal/Ocular Effects. Burns and irritation of the skin and eyes are the most frequent manifestations of coal-tar creosote toxicity following

TABLE 2-2. Levels of Significant Exposure to Creosote - Dermal

Figure	•	Exposure Frequency/			LO		٠		
Key	Species	Duration	Effect	NOAEL	Less Serious		Serious	Reference	Form
CUTE EXP	OSURE		•						
Death	• .								
1	Rabbit	1 d 1x/d				> 7.95	g/kg (LD50)	Pfitzer and Gross 1964	coal
Systemic						•			
2	Mouse	7-21 d 7d/wk 1x/d	Derm	pa	5% in (irritation)			Wrench and Britten 1975	coal
NTERMEDIA	TE EXPOSURE					•			
Cancer				•					
3	Mouse	8 mo 2x/wk	Derm			0.25ml	(CEL)	Ros et al. 1958	coal
4	Mouse	44 wk 2x/wk	Derm			0.25ml	(CEL)	Boutwell and Boach 1958	coal
HRONIC EX	POSURE								
Cancer				•					
5	Mouse	70 wk 2x/wk	Derm	·		1 drop undiluted	(CEL)	Lijinsky et al. 1957	coal
6	Mouse	Life 3x/wk	Derm				(CEL)	Poel and Kammer 1957	coal

d = day(s); wk = week(s); x/d = times per day; x/wk = times per week; Derm = dermal; Resp = respiratory; CEL = cancer effect level; ml = milliliters; > = greater than.

dermal exposure. Creosote burns were observed in construction workers who handled wood treated with creosote (presumably coal-tar creosote) (Jonas 1943). Exposure levels were not specified. It was found that 70% of the burn cases were mild and were characterized by erythema of the face. These symptoms were more marked on the cheeks, nose, forehead, and posterior part of the neck. The remainder of the burn cases (30%) were more severe and were characterized by intense burning, itching, and considerable subsequent pigmentation followed by desquamation. There is no way to determine, given the information provided, whether this response was due to primary irritation or an allergenic response. Conjunctivitis was seen in 15% of the men, with 3% of these showing corneal injuries. Exposure to the sun exacerbated these effects. The skin injuries were reversed with thorough washing, application of Foille solution (which contains various calcium salts, benzocaine, sulphur and vegetable oil base), and cessation of exposure.

Other studies confirm that coal-tar creosote is capable of inducing phototoxicity of the skin. Phototoxicity is an exaggerated response to sun exposure characterized by excessive sunburn. The usual sunburn is produced by UVB light, whereas a phototoxic skin response occurs following exposure to UVA light (Baker and Fannick 1981). Coal-tar creosote exposure was evaluated in six dock builders, all male of Scandinavian descent with fair features and an average of 16.6 years spent as pile-drivers (Baker and Fannick 1981). It was found that the dermal and ocular burning and irritation experienced by these workers upon dermal contact with wood treated with coal-tar creosote was exacerbated on hot or sunny days. Skin examinations of these dermally exposed workers revealed erythema and dry peeling skin on the face and neck with irritation and folliculitis on the forearms. These symptoms were worse on hot or sunny days, at which time red, swollen, and puffy eyes were observed. Thus, phototoxicity compounds the irritative response of the skin to coal-tar creosote. No skin tumors were observed. Exposure levels of coal-tar-pitch volatiles measured were considered to represent the minimum ambient air exposure in the barge area because of unusual environmental sampling conditions. These levels ranged from below the detectable limit to 0.06 mg/m<sup>3</sup> for breathing zone samples and from below the detectable limit to 0.02 mg/m<sup>3</sup> for area samples.

Coal-tar creosote has been reported to produce other types of noncancerous skin lesions as well as burns and irritation following dermal exposure. Haldin-Davis (1935) described the case of a man employed in the activity of dipping wood in creosote tanks who received "heavy" dermal exposure to creosote (level not determined) on the face, trunk, and thighs. He subsequently developed a number of lesions on the hands, forearms, and thighs. One of these lesions was excised and examined and classified as a (benign) squamous cell papilloma.

The National Institute for Occupational Safety and Health (NIOSH) was called in to investigate potential employee-related health effects in mixers and laborers exposed to coal-tar products in a factory (Chrostek 1980). Six

of seven environmental samples collected exceeded the NIOSH Permissible Exposure Limit (PEL) for coal-tar products (cyclohexane solubles, which includes creosote) of 0.1 mg/m³. Exposures were reported to be as high as 1.42 mg/m³. Because environmental levels significantly exceeded the PEL, it is likely that considerable inhalation exposure occurred. One worker reported the need for medical treatment when creosote splashed in his eye, and another claimed that he was hospitalized for 3 days after experiencing convulsions following an incident where he was splashed with creosote. The medical records for these two men were not reviewed.

Skin irritation is also observed in laboratory animals following dermal exposure to creosote. Rabbits given single dermal applications of undiluted coal-tar creosote exhibited slight to moderate erythema and moderate edema followed by severe hyperkeratosis (Pfitzer and Gross 1964). These same investigators also found that instillation of 0.1 mL undiluted coal-tar creosote in the eyes of rabbits produced a redness of the conjunctiva with congested vessels that resolved within seven days (Pfitzer and Gross 1964). The effects of dermally applied coal-tar fractions, derived from creosote and anthracene oils by high-temperature boiling, were studied in mouse tail skin at concentrations of 5% and 10% (acids) in paraffin by Wrench and Britten (1975). Several of the fractions caused irritation, and some caused peeling and epidermal thickening. The authors concluded that the acids that boiled in the range from 280° to 340°C have a more specific action in inducing granular layers than the parent tars, oils, or whole acids at similar concentrations. This study is of limited value because the chemical composition of the acid fractions was not defined, and no dose levels could be quantified. It can only be concluded that certain fractions of coal-tar creosote irritate mouse skin.

Creosote also irritates other types of tissue. The effects of beechwood creosote (dose not specified) on the periapical tissue (the connective tissue surrounding the apex of the tooth) were studied in 10 teeth from 4 different dogs 7 days after its application following root canal surgery (Attalla 1968). Beechwood creosote application resulted in localized inflammatory changes and occasional abscess formation in the periapical region, presumably due to tissue damage caused by coagulation of proteins. Bone resorption was also observed in the alveolar process. The authors concluded that the beneficial disinfectant properties of creosote may be outweighed by its irritant effect following root canal surgery.

In summary, dermal application of either coal-tar or beechwood creosote results in mild-to-severe irritation of the skin, connective tissue, and eyes, as well as benign skin lesions in both humans and animals. Coal-tar creosote also induces phototoxicity, so exposure to the sun exacerbates its irritant effects. This has important implications for individuals living in areas surrounding hazardous waste sites who may come in contact with creosotecontaminated soils.

# 9 3 3 Immunological Effects

Several cases of acute allergic dermatitis have been reported following contact with creosote bush resin. Smith (1937) described the case of a patient who presented with erythematous and vesicular dermatitis of the face, the upper part of the neck, and the backs of the hands after collecting creosote bush. Patch tests confirmed the existence of an allergy to this plant. Leonforte (1986) reported six cases of acute allergic dermatitis subsequent to contact with a creosote bush and confirmed by a patch test. Two cases were a result of "casual occupations," two were a result of household remedies, and two were a result of burning the bush. Based on his findings, the author concluded that the allergens are probably contained in the plant's perfume, are volatile, and are not destroyed by heat. The relevance of these findings to individuals who live in areas surrounding hazardous waste sites and who will most likely be exposed to coal-tar creosote is questionable. Creosote bush resin differs from creosote extracted from coal and wood tar. but all contain phenolic derivatives. It is not known whether these derivatives are the allergens in creosote bush resin.

# 2.2.3.4 Neurological Effects

A worker who was splashed with creosote in a refinery claimed that he was hospitalized for 3 days after experiencing convulsions (Chrostek 1980). The medical records for this man were not reviewed, so it cannot be concluded that creosote exposure was responsible for his convulsions.

No studies were located regarding neurologic effects in animals following dermal exposure to creosote.

### 2.2.3.5 Developmental Effects

No studies were located in the literature regarding developmental effects in humans following dermal exposure to creosote.

The only available study in animals reported that dermal contact with creosote-treated wood produced fetotoxic effects in pregnant sows (Schipper 1961). Four sows were confined to wooden farrowing crates for 2-10 days before delivery. The platforms of the crates were coated with 3 brush applications of a commercial wood preservative containing 98.5% coal-tar creosote. Following this gestational exposure to creosote, 24 of the 41 pigs delivered were dead at birth, and 11 pigs died by day 3 post-farrowing. The surviving pigs had rough skin and suffered from dehydration and severe diarrhea. The pigs failed to gain weight until they were 5-6 weeks old. No toxic effects on the sows were reported. On the contrary, four sows confined to untreated lumber crates at least 24 hours before farrowing delivered 36 pigs; 1 died within 24 hours and 3 died post-farrowing. No toxic effects were noted in mothers or baby pigs. The study limitations include unequal duration of exposure between treated and untreated groups, and the lack of

statistical analysis of the reported results. Moreover, although creosote was absorbed in sufficient quantities via dermal contact by pregnant sows, the actual dose producing toxic effects in pigs cannot be determined. Therefore, neither dose-response nor NOAEL and LOAEL values for the developmental toxicity of creosote for pigs can be determined.

# 2.2.3.6 Reproductive Effects

No studies were located regarding reproductive effects in humans or animals following dermal exposure to creosote.

#### 2.2.3.7 Genotoxic Effects

No studies were located regarding genotoxic effects in humans or animals following dermal exposure to creosote. The mutagenic activity of urine samples taken from workers briefly exposed to coal-tar creosote vapors in a wood preserving factory was investigated (Bos et al. 1984a, see Section 2.2.1.7, "Inhalation Exposure, Genotoxic Effects"). However, the possibility of dermal exposure by contact with residual surface creosote cannot be excluded in this study. The urine samples were negative for mutagenic activity following exposure at work. The data are not sufficient to draw a conclusion regarding the genotoxic potential of dermal contact with creosote vapors in humans, but they do suggest that extrapolation of mutagenicity tests on a single component (i.e., PAHs) may not predict the action of the mixture.

# 2.2.3.8 Cancer

Various case reports and the results of cross-sectional occupational surveys associate chronic occupational creosote exposure with the development of skin cancer (Cookson 1924; Henry 1947; Lenson 1956; MacKenzie 1898; O'Donovan 1920). These papers reported a similar disease etiology for different groups of workers exposed to creosote that included the development of dermatoses, such as squamous papillomas, that progressed to carcinoma, usually squamous-cell carcinoma. The latency period for the development of dermatoses, such as squamous papillomas, was usually 20-25 years (Cookson 1924; Henry 1946, 1947; O'Donovan 1920). The latency period for the development of carcinoma averaged 25 to 30 years (Cookson 1924; Henry 1946, 1947; O'Donovan 1920). Worker exposure in the past was much greater than it now is because of less-sophisticated industrial practices used in the past, the lack of knowledge concerning occupational hygiene, and the current recognition of the dangers of excessive exposure to the health of workers. Other factors that should be considered when extrapolating the findings of this older literature to present conditions are the role of exposure to ultraviolet radiation in the form of sunshine in these workers (ultraviolet radiation is now known to be a major cause of skin cancer), the composition of the creosote products, and other health factors that differ in Great Britain prior to 1940 and the present. Although these studies lack information concerning specific exposures and did not consider other risk factors for the

development of skin cancer, when taken as a group, they present convincing evidence of a relationship between chronic creosote exposure, phototoxicity, and the development of skin carcinoma in humans.

Creosote exposure has also been associated with multiple myeloma in a case/referent study of a Swedish population (Flodin et al. 1987, see Section 2.2.1.8). This relationship needs to be investigated further by more sophisticated studies.

On the other hand, several studies are available that suggest that there is no association between exposure to creosote or other coal-tar products and cancer in humans. No increased incidence of non-melanoma skin cancer was observed in 426 patients 25 years after they had undergone 4 years of coal-tar medicinal therapy in combination with ultraviolet light for the treatment of atopic dermatitis and neurodermatitis (Maugham et al. 1980). This study is limited in that follow-up occurred in only 72% of the patients, and there was no discussion of recall bias or of the effects of mobility (i.e., relocation of the study participants) on the results. In a 25-year retrospective study of 280 psoriatic patients treated with crude coal-tar in combination with ultraviolet radiation at the Mayo clinic, it was found that the incidence of skin cancer in these patients was not significantly increased above the expected incidence for the general population (Pittelkow et al. 1981). However, it should be noted that in investigations on populations using therapeutic coal-tar skin products, exposure was to a different coal tar mixture. Creosote is a lower boiling distillate fraction of crude coal-tar than the components contained in the skin products, and creosote does not contain the concentrations of putative carcinogenic PAH compounds that are present in the coal-tars contained in the skin products.

Another study was located that reported that there was no increase in the risk of skin, bladder, or lung cancer in wood treatment plant workers (Tabershaw Occupational Medicine Associates 1980). Limitations in this study include: the study population was small; the study population was comprised of 46.5% blacks who experience a very low incidence of skin cancer as compared to whites, thus biasing the results; the exposure and follow-up periods did not allow a long enough latency period for tumor development; and there was no verification provided that those studied were actually exposed to creosote or coal-tar.

A large body of evidence exists that coal-tar creosote is carcinogenic when applied to the skin of laboratory animals. Many of the early studies are limited in that they lack appropriate negative control data, the dose of creosote and the chemical composition of the fractions studied were not quantified, and no other tissues were generally examined. The results from later studies that include appropriate control groups are consistent with the earlier studies that found that creosote is carcinogenic following dermal application to rodent skin. Six representative dermal carcinogenicity studies

are reviewed below, and the relevant cancer effect levels are presented in Table 2-2.

The tumor-promoting potential of creosote was evaluated by applying various fractions (e.g., basic, phenolic, and neutral) to the skin of mice in conjunction with benzo(a)pyrene (B[a]P) (Cabot et al. 1940). The various fractions of creosote oil were prepared by distillation and separation. Ninety percent of the creosote oil distilled between  $160^\circ$  and  $300^\circ$ C. The basic fraction was removed with aqueous hydrochloric acid. The phenolic fraction was removed with aqueous sodium hydroxide. The remaining neutral fraction was then steam distilled. The composition of the various fractions was not specified. Four of the test solutions antagonized the tumorigenic effects of B[a]P, but in three instances this antagonistic effect was considered to be secondary to skin damage. Only the phenolic fraction seemed to exhibit a primary antagonistic effect. Three fractions (basic, neutral, and neutral distillate) exhibited apparent promoting effects.

The potential of basic fractions of creosote to accelerate tumor induction by known carcinogens was evaluated by Sall and Shear (1940). A 1% solution of the basic fraction of creosote oil in benzene was dermally applied to female strain A mice alone or in conjunction with 0.05% or 0.02% B[a]P. The basic fraction alone did not induce skin tumors, but when applied in conjunction with either concentration of B[a]P, skin tumors appeared more rapidly than when B[a]P alone was applied. Maximum tumor induction was seen between 28 and 42 weeks; 19 or 20 mice developed tumors. The composition of the basic fraction was not specified, but it is reported to be noncarcinogenic.

The ability of creosote to induce lung tumors after dermal application to mice was studied when it was observed that mice housed in creosote-treated wooden cages had a high incidence of lung tumors (Roe et al. 1958). Dermally applied creosote (0.25 mL undiluted twice weekly for 8 months) induced 5.8 lung adenomas/mouse in mice that were reared in stainless steel cages. Creosote treatment following the same regimen in mice reared in creosotetreated cages induced 10.8 lung adenomas/mouse. Untreated controls reared in untreated cages exhibited 0.5 lung adenomas/mouse. A high incidence of skin tumors was also observed in the creosote-treated mice reared in either type of cage. In a second experiment, topical application of "one drop" of creosote twice a week for only 4 weeks induced lung adenomas but not skin tumors in mice reared in stainless steel cages. This study demonstrated that creosote induces tumors in the lungs and skin of mice when dermally applied. Rearing animals in creosote-treated wooden cages exacerbated the tumorigenic effect of dermally applied creosote. Based on this study, lung tumors may be a more sensitive end point of creosote tumorigenic activity than skin tumors.

Seven groups of thirty female albino mice each were treated with dimethylbenzanthracene (DMBA) and creosote (0.25 mL of undiluted creosote oil), alone and combinations of the two to evaluate the carcinogenic,

Initiating, and promoting activity of creosote on mouse skin (Boutwell and Bosch 1958). Creosote alone, DMBA and creosote, and creosote with croton oil all induced the development of papillomas and carcinomas. Tumors first appeared at 10-20 weeks of application. DMBA pretreatment shortened the latent period and increased the tumor yield of creosote treatment, but since almost a nearly maximal tumor induction response (i.e., percent of mice with papillomas was just below 100) was seen with creosote alone, tumor-promoting activity of creosote could not be definitely proven. The initiating activity of creosote was demonstrated by its ability to induce tumors when applied prior to croton oil treatment (croton oil alone was without effect). Creosote alone or in combination with DMBA or croton oil induced papilloma formation more slowly, and carcinomas appeared by 14 weeks and accumulated more rapidly than in the DMBA plus croton oil group. Thus, this study demonstrated the carcinogenic and tumor-initiating activity of creosote on mouse skin.

The complete carcinogenic and tumor-promoting activity of undiluted creosote, a 10% solution of creosote oil, and 2% solution of the basic fraction of creosote was studied when dermally applied to mice (Lijinsky et al. 1957). Undiluted creosote alone induced 23 skin tumors (16 malignant) in 13 of 26 treated mice with a latent period of 50 weeks. The authors concluded that creosote alone has a carcinogenic activity comparable to a 0.01% solution of dimethylbenzanthracene (DMBA). When applied as a promoter following a single application of DMBA, 32 skin tumors (26 malignant) were observed in 17 of 30 mice with a latent period of 39 weeks. The basic fraction did not act as a tumor promoter when administered after a single application of DMBA. Thus, creosote appeared to enhance the carcinogenic activity of DMBA, but the promoting effect was not strong, when the results were compared to DMBA positive controls.

The carcinogenic activity of two high-temperature-derived creosote oils ("light" and "blended") was studied by Poel and Kammer (1957). The principal components of light oil are benzene, toluene, xylene and solvent naphtha. Blended oil is a mixture of creosote oil, anthracene oil and the oil drained from the naphthalene recovery operation. Its principal components are methylated naphthalenes, acenaphthene, fluorene, phenanthrene, anthracene and carbazole. The oils were applied by drops to the skin of mice at concentrations of 20%, 50%, or 80% three times a week for life. Both oils induced skin tumors in every exposed mouse by 21 to 26 weeks of application. Several mice exhibited metastases to the lungs or regional lymph nodes. The fractions tested did not contain B[a]P, so the authors concluded that the carcinogenic activity of creosote was not due to B[a]P.

In conclusion, the results of these studies indicate that coal-tar creosote and several of its fractions are carcinogenic when applied to the skin of mice. Dermally applied creosote can also act as a tumor-initiating agent when applied prior to croton oil treatment and can enhance and accelerate tumor induction by B[a]P. Thus, it is likely that individuals whose skin comes into contact chronically with creosote would be at higher

risk for cancer, particularly when exposure to other carcinogenic substances also occurs, as is a likely scenario in areas surrounding hazardous waste sites.

#### 2.3 TOXICOKINETICS

Specific information regarding the toxicokinetics of creosote is limited. Several compounds have been detected in creosote, yet there are no definitive data on which of these compounds people are exposed to in wood-treatment plants or at hazardous waste sites. Analyses have revealed that PAHs are the major components of the creosote mixture (EPA 1981a). Hence, pharmacokinetic studies on PAHs can be used as surrogates for creosote. However, this is only speculative given the possible toxicokinetic interactions that occur among the PAHs and other components in the creosote mixture, and will be used only when data on creosote are not available. For more information on the toxicokinetics of PAHs, please refer to the ATSDR Toxicological Profile for Polycyclic Aromatic Hydrocarbons (ATSDR 1990).

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## 2.3.1 Absorption

### 2.3.1.1 Inhalation Exposure

No studies were located in humans or animals regarding the direct analysis of the extent or rate of creosote absorption following inhalation exposure. However, there is evidence to suggest that inhalation absorption occurs. Employees of a creosote-impregnating plant excreted 1-hydroxypyrene, a metabolite of the creosote component, pyrene, in their urine (Jongeneelen et al. 1985). The presence of this metabolite in the urine suggested that creosote components were absorbed and metabolized following inhalation exposure. However, it is also possible that some dermal exposure may have occurred as well.

PAHs extracted from coal fly ash were intratracheally administered to pregnant rats at a dose of 20 mg/kg. The presence of the PAHs in the maternal and fetal lungs and livers indicated that pulmonary absorption occurred (Srivastava et al. 1986).

### 2.3.1.2 Oral Exposure

No studies were located in humans and animals regarding the direct analysis of the extent or rate of creosote absorption following oral exposure. However, the presence of creosote metabolites in the urine of humans and rabbits receiving calcium creosotate (a calcium salt of creosote) tablets was evidence that creosote was absorbed following ingestion (Fellows 1937, 1939b). Furthermore, evidence exists that certain PAHs found in creosote such as anthracene (Rahman et al. 1986), benzo(a)pyrene (Hecht et al. 1979; Rahman et al. 1986; Rees et al. 1971; Yamazaki et al. 1987), chrysene (Chang 1943;

Modica et al. 1983), and phenanthrene (Rahman et al. 1986) are absorbed following oral administration in animals.

# 2.3.1.3 Dermal Exposure

No studies were located in humans and animals regarding the direct analysis of the extent or rate of creosote absorption following dermal exposure. However, reports of workers who developed cancer subsequent to dermal exposure suggested that creosote was absorbed through the skin (Gookson 1924; Henry 1946, 1947; Lenson 1956). It can also be concluded that dermal absorption occurred as evidenced by the development of skin tumors (Boutwell and Bosch 1958; Lijinsky et al. 1957; Poel and Kammer 1957; Roe et al. 1958) and lung tumors (Roe et al. 1958) in mice following the dermal application of creosote. Based on these occupational exposures and animal toxicity studies, it would appear that dermal absorption of creosote occurs.

# 2.3.2 Distribution

## 2.3.2.1 Inhalation Exposure

No studies were located in humans or animals regarding the distribution of the creosote mixture following inhalation exposure. However, an occupational study indicated that creosote is absorbed and metabolized as evidenced by the appearance of 1-hydroxypyrene (a metabolite of pyrene, a component of creosote) following inhalation exposure (Jongeneelen et al. 1985). Because creosote is composed of hydrocarbons, creosote is likely to distribute to lipid-rich tissues. Creosote is also likely to distribute to the liver as evidenced by the presence of metabolites in the urine.

When [3H]-benzo[a]pyrene was administered intratracheally to rats at a dose of 0.001 mg/kg, radioactivity was distributed to all tissues. During the 6 hours following administration, more than 20% of the dose was detected in the carcass. The activity steadily increased in the intestine and the intestinal contents over the 6 hours. Levels of activity in the liver and lung were moderate and declined over time. Trace amounts of activity were detected in other tissues (Weyand and Bevan 1987).

Intratracheal administration of [3H]-benzo[a]pyrene along with the benzene extract of coal fly ash to pregnant rats (20 mg/kg/day) on days 18 and 19 of gestation resulted in their distribution to the maternal lung and liver. The amount of radioactivity found in the maternal liver was approximately 68% of the amount of radioactivity found in the maternal lung. The amount of radioactivity found in the placenta, fetal lung, and fetal liver was approximately 4%, 1.9%, and 1.4%, respectively, of the amount of radioactivity found in the maternal lung (Srivastava et al. 1986). Much of the radioactivity was attributable to metabolites. These results suggest that components of creosote and their metabolites can pass through the placenta and distribute to fetal tissue.

## 2.3.2.2 Oral Exposure

No studies were located in humans or animals regarding the distribution of creosote following ingestion. Based on chemical structure, it is likely that PAHs would have a strong affinity for adipose tissue. Benz[a]anthracene, chrysene, and triphenylene distributed to all tissues following oral administration (22.8 mg/kg) to female rats, but its greatest distribution was to adipose tissue. In this study, benz[a]anthracene concentrations in adipose were 10 times higher than in other tissues (Bartosek et al. 1984).

The distribution of nonmetabolized PAHs is dependent on their water-solubility. The more water-soluble PAHs, like triphenylene, are generally more available to tissues other than fat (Bartosek et al. 1984). In humans, distribution of creosote following ingestion is likely to be qualitatively similar to that seen in the animal studies. The lipophilicity of PAHs allows the chemicals to be readily absorbed and preferentially accumulated in fatty tissues. Furthermore, PAHs are likely to be present in adipose and highly perfused organs such as the lungs and liver.

# 2.3.2.3 Dermal Exposure

No studies were located in humans or animals regarding the distribution of creosote following dermal exposure. Distribution of creosote in humans following dermal exposure is expected to be qualitatively similar to that seen in animals or in humans following any route of exposure.

### 2.3.3 Metabolism

Generally, the PAH components of creosote are metabolized by oxidative enzymes to generate active metabolites that can bind to macromolecules. The metabolic profiles vary among species and compounds, but the components follow the same reaction pathways. Hence, the metabolites are structurally very similar. The proposed metabolic scheme for a representative PAH, benzo(a)pyrene is presented in Figure 2-2. The principal products include phenols, dihydrodiols, quinones, anhydrides and conjugates of these products (Autrup and Seremet 1986; Dahl et al. 1985; Geddie et al. 1987; Hopkins et al. 1962; Jongeneelen et al. 1985; Petridou-Fischer et al. 1988; Povey et al. 1987; Rice et al. 1986; Weyand and Bevan 1987).

# 2.3.3.1 Inhalation Exposure

Workers in a creosote-impregnating plant were exposed to creosote by inhalation. The creosote that these employees inhaled contained 19.8 mg pyrene/g creosote (approximately 2%). A metabolite of pyrene, 1-hydroxypyrene, was detected in their urine at levels that were above the mean values of controls (Jongeneelen et al. 1985). The identification of PAH metabolites in the urine could serve as a method of biological monitoring of

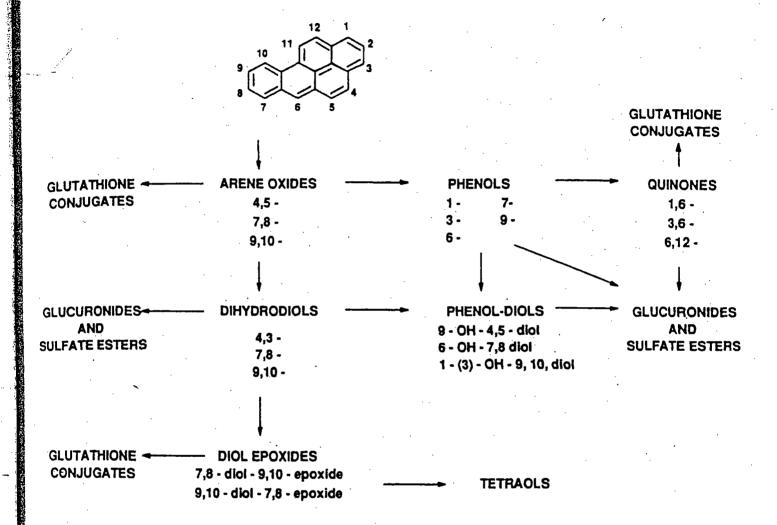


FIGURE 2-2. Proposed Metabolic Scheme for Benzo(a)pyrene

exposed workers, and possibly individuals living in the vicinity of hazardous waste sites where creosote has been detected.

## 2.3.3.2 Oral Exposure

Calcium creosotate was orally administered to humans at daily doses of 7-30 mg/kg for 3 days. Calcium creosotate phenols were excreted in the urine. Also, large unspecified doses of calcium creosotate were orally administered to rabbits. Analysis of the rabbit urine revealed that free and conjugated phenols were excreted (Fellows 1939b).

Rats receiving a single dose of either 0.0002, 0.002, 0.02, 0.2, or 2.0 mg pyrene/kg by gavage in olive oil excreted 1-hydroxypyrene in the urine in a dose-dependent manner. This metabolite could be detected up 96 hours after administration. No pyrene was excreted (Jongeneelen et al. 1985).

It is evident in both human and animal studies that hydroxylation is a principal oxidative pathway of PAH metabolism, and consequently, creosote metabolism. In these studies, there were no discussions to suggest that the researchers attempted to identify other metabolites.

# 2.3.3.3 Dermal Exposure

Two patients suffering from eczema on the arms and legs were treated for several days with an ointment containing 10% pix lithanthracis dermata (coaltar). The daily dermal dose was approximately 1 mg/kg. Analysis of the urine samples collected from these patients prior to treatment and in the morning and evening of the first three days of treatment showed that 1-hydroxypyrene (a metabolite of pyrene which is a PAH component of coal-tar) was excreted at levels 200 times that which was detected before the treatment started (Jongeneelen et al. 1985). Thus, based on these results, it appears that PAH components of coal-tar creosote are metabolized following dermal exposure in humans.

### 2.3.4 Excretion

Excretion of the PAH compounds of creosote is controlled by their rate of metabolism. Excretion of these metabolites or any remaining parent compound is primarily in the bile and the feces. Weyand and Bevan (1987) demonstrated this by cannulating the bile ducts of rats that received [<sup>3</sup>H]-benzo[a]pyrene intratracheally. Those rats with the biliary cannulas had significantly lower levels of activity in their intestines, intestinal contents and stomach than rats without biliary cannulas. Sanders et al. (1986) showed that PAHs were primarily removed in the feces after dermal administration of [<sup>14</sup>C]-benzo[a]pyrene and [<sup>14</sup>C]-7,12-dimethylbenz[a]-anthracene, suggesting hepatobiliary excretion. Urinary excretion also occurs, but to a lesser extent than the other routes.

# 2.3.4.1 Inhalation Exposure

No studies were located in humans or animals regarding the excretion of creosote following inhalation exposure.

## 2.3.4.2 Oral Exposure

No studies were located in humans or animals regarding the excretion of creosote following oral exposure.

## 2.3.4.3 Dermal Exposure

No studies were located in humans or animals regarding the excretion of creosote following dermal exposure.

#### 2.4 RELEVANCE TO PUBLIC HEALTH

Creosote is a complex mixture of organic substances. It can be derived by distillation from coal (coal-tar creosote) or wood (beechwood creosote). Another form is a resin from the leaves of the creosote bush. The chemical composition of creosote varies considerably depending on the source of the coal, wood, or plant and the design and attendant operating conditions (temperature, gas distillation systems, etc.) used to produce the creosote. Coal-tar creosote is reported to contain over 300 compounds (EPA 1978), the major components of which are PAHs, tar acids (phenol, cresols and xylenols), and tar bases (pyridine and lutidine derivatives). Coal-tar creosote is used primarily as a wood preservative. Beechwood creosote consists mainly of phenol, cresol, guaiacol, xylenol and creosol. It has been used therapeutically as an expectorant and a disinfectant. Creosote bush resin consists of phenolics (which account for 83-91% of the total resin), neutrals (e.g., waxes), basics (e.g., alkaloids), and acidics (e.g., phenolic acids). Given the current widespread use of coal-tar creosote as a wood preservative and its past pesticidal applications, it is the form of creosote most likely to be present at hazardous waste sites.

There is relatively little information available regarding the systemic effects of ingested creosote. However, oral exposure via ingestion of contaminated drinking water is most likely not a highly significant route of exposure to creosote at hazardous waste sites, given its low solubility in water. Most of the available recent information on the adverse effects of ingested creosote in both humans and animals is on beechwood creosote. These studies indicate that the liver, kidney, central nervous system, and cardiovascular system may be adversely affected by beechwood creosote following primarily acute exposure, although the mechanism of action for these effects is not known. The relevance of these findings in animals with exposure to beechwood creosote to humans exposed to coal-tar creosote is not known. Although the two forms of creosote contain some components in common

(e.g., phenolic derivatives) it is not known whether the toxic effects observed after exposure to beechwood creosote can be attributed to the phenol components and thus, extrapolated to coal-tar creosote. Coal-tar creosote exerts its acute toxic effects primarily via dermal exposure, causing architectural damage to the tissues that it comes in contact with, such as the skin and eyes. Acute ingestion of coal-tar creosote appears to affect primarily the kidney and liver. The mechanism of action for coal-tar creosote-induced toxicity is not defined, but is most likely due to the activity of the PAH components.

Much of the information available on the health effects of coal-tar creosote must be inferred from experimental animal and <u>in vitro</u> data on the components of the creosote mixture (i.e., PAHs). Extrapolation of these results to possible human health effects following exposure to creosote must consider the possible interactions of the mixture, such as co-carcinogenicity, co-mutagenicity, additivity, promotion, and antagonism. The extent to which these interactions modify the expression of creosote toxicity in humans is not known.

Death. No information was available on the lethal effects of creosote following inhalation or dermal exposure in humans. No deaths were observed in rats exposed to near saturated vapors of coal-tar creosote for one hour (concentration estimated to be less than 0.033 mL/L) (Pfitzer and Gross 1964). The exposed animals showed signs of slight eye and nose irritation and slight dyspnea. All rats exhibited weight gains comparable to the control animals and there were no treatment-related lesions observed at necropsy. The dermal LD $_{50}$  in rabbits has been estimated to be greater than 7.95 g/kg (Pfitzer and Gross 1964).

Ingestion of creosote can be fatal to both humans and animals. case reported by Bowman et al. (1984), death following ingestion of creosote (form and quantity not specified) was attributed to multi-organ failure. Death has been reported to occur in adults and children 14 to 36 hours after the ingestion of about 7 g or 1-2 g creosote, respectively (Clayton and Clayton 1981). The latter anecdotal information is of limited value because it was obtained from a secondary source that provided no supporting documentation. The oral  $LD_{50}$  for coal-tar creosote is reported to be 433 mg/kg in the mouse and 725 mg/kg in the rat (RTECS 1988). However, another study reported an acute oral  $LD_{50}$  of 1700 mg/kg in the rat (Pfitzer and Gross 1964). The reasons for this discrepancy in  $LD_{50}$  values are not known. Based on these data, coal-tar creosote can be classified as mildly to moderately toxic. Beechwood creosote is lethal to animals, although the doses required to produce death are relatively high (300-525 mg/kg). Death is preceded by signs of central nervous system intoxication following acute exposure in both humans and animals. Given the relatively high oral doses required to cause death following acute exposure in both humans and animals, individuals living

in areas surrounding hazardous waste sites contaminated with creosote may not be at high risk for death due to acute ingestion of creosote.

Systemic Effects. A slightly higher incidence of bronchitis or thickening of the tracheal mucous membrane was observed in mice fed beechwood creosote for 52 weeks. This effect was considered secondary to irritation resulting from long-term inhalation exposure to volatile (but unidentified) components of creosote in the feed. No toxic respiratory effects have been reported following coal-tar creosote exposure in either humans or animals. The airborne component chemicals in coal-tar creosote have not been well defined. However, based on the results discussed above, it can be assumed that some components of beechwood creosote, that may be common to coal-tar creosote are irritating to the respiratory tract.

A report was found in the older literature that described the case of a woman who experienced hypertension attributed to creosote exposure by "self-medication" for chronic bronchitis (Robinson 1938). Cardiovascular collapse has also been reported to occur following the ingestion of lethal doses of coal-tar creosote in humans (Clayton and Clayton 1981). These cases are anecdotal and occur in isolated instances. There is often a lack of exposure data, and the possibility of confounding factors limit the usefulness of these findings. Furthermore, adverse creosote-induced cardiovascular effects have not been reported in animals. Thus, it is not likely that the cardiovascular system is a major target organ of toxicity from creosote exposure.

Ulceration of the oropharynx and petechial hemorrhages over the gastrointestinal serosal surfaces were noted at autopsy in the case of a 70-year-old man who died following the ingestion of "industrial" creosote (Bowman et al. 1984). This acute tissue damage was attributed to a corrosive action of phenol (a component of creosote) via its ability to denature and precipitate proteins. Animals that died following the administration of single gavage doses of coal-tar creosote in an acute range-finding study (doses ranged from 2.52-5.00 mg/kg) exhibited hyperemia and distention of the stomach upon necropsy. Similarly, rabbits that died following single dermal applications of undiluted coal-tar creosote exhibited hyperemia of the intestines (Pfitzer and Gross 1964). The toxicological significance of these changes is not known. Thus, given that gastrointestinal lesions were observed in only one isolated instance in humans, and the fact that only mild creosoteinduced gastrointestinal effects of unknown toxicological significance have been reported in animals, it is likely that the gastrointestinal system is a target organ of toxicity only after the ingestion of relatively high doses of creosote.

Various hematological and clinical chemistry parameters have been observed to be altered by dietary exposure to beechwood creosote in rats and mice (Miyazato et al. 1981, 1984a,b). The only effect considered to be treatment-related by the authors was an increase in serum cholesterol in both

rats and mice. The toxicological significance of this change is not known, particularly with regard to individuals exposed to coal-tar creosote in areas surrounding hazardous waste sites. No adverse hematological effects have been reported in humans exposed to creosote.

Degeneration and necrosis of hepatocytes were observed at autopsy in the case reported by Bowman et al. (1984). Given the advanced age of this man, the possibility of confounding factors, and the lack of comparison data, it is not possible to definitively attribute these effects to creosote ingestion. Dietary exposure to beechwood creosote induces changes suggestive of liver injury in rats and mice (increase relative liver weights and increased liver enzymes). However, no treatment-related changes were observed at histological evaluation (Miyazato et al. 1981, 1984a,b), and the toxicological significance of these changes with regard to individuals exposed to coal-tar creosote in areas surrounding hazardous waste sites is not known.

Observations made in both humans and animals suggest that creosote has the potential to induce adverse effects in the kidney. Prior to death, the patient described by Bowman et al. (1984) became acidotic and anuric, indicating kidney failure. Acute renal tubular necrosis was revealed at necropsy. However, the acute tubular necrosis may have been due to vascular insufficiency rather than a direct toxic effect on the kidney. Male rats chronically exposed to beechwood creosote in the diet exhibited an exacerbation of the spontaneously occurring chronic nephrosis normally seen in aging male rats (Miyazato et al. 1984b). Thus, the kidney may be adversely affected by both acute high-level exposure (e.g., following an accidental spill) to coal-tar creosote and chronic low-level exposure to beechwood creosote.

Burns, irritation and benign lesions of the skin (e.g., squamous papillomas), and conjunctivitis of the eyes are the most frequent manifestations of coal-tar creosote toxicity following dermal exposure in both humans and animals. These effects consist of mild-to-severe erythema, intense burning, itching, and subsequent pigmentation followed by desquamation. Involvement of the skin, connective tissue, and eye are often seen. Creosote is also capable of inducing phototoxicity of the skin, so exposure to sun exacerbates its irritant effects. This has important implications for individuals living in areas surrounding hazardous waste sites who may come in contact with creosote-contaminated soils.

Immunological Effects. No adverse immunologic effects have been reported in either humans or animals following exposure to coal-tar or beechwood creosote. Several cases of acute allergic dermatitis have been reported following contact with creosote bush resin. The relevance of these findings to individuals who live in areas surrounding hazardous waste sites and will most likely be exposed to coal-tar creosote is not known. Creosote bush resin differs from creosote extracted from coal and wood tar, but all

contain phenolic derivatives. It is not known whether these derivatives are the allergens in creosote bush resin.

Neurological Effects. Adverse neurologic effects have been reported in both humans and animals following the acute ingestion of high doses of coaltar or beechwood creosote. Such effects include salivation, vomiting, respiratory distress, thready pulse, vertigo, headache, loss of pupillary reflexes, hypothermia, cyanosis, muscle twitching, and convulsions. These observations suggest that creosote may be a general central nervous system stimulant following acute high-level exposure. However, the possibility exists that some of the effects observed in animals that licked treated utility poles may have been due to ingestion of pentachlorophenol. Pentachlorophenol, like creosote, is an oil-borne wood preservative with extensive use in the public utility industry for treatment of utility poles. The neurotoxic effects observed in cattle that licked treated utility poles are more compatible with the metabolic effects (i.e., uncoupling of phosphorylative oxidation) associated with pentachlorophenol (for more information on the effects of pentachlorophenol, please refer to the ATSDR Toxicological Profile for Pentachlorophenol, ATSDR 1989a). Thus, it is not possible to determine conclusively whether creosote is toxic to the central nervous system of animals based on this report. However, no information was found to suggest that chronic low-level exposure to creosote by individuals in areas surrounding hazardous waste sites would result in neurotoxicity.

Developmental and Reproductive Effects. The only available study in animals reported that dermal exposure to creosote-treated wood produced fetotoxic effects in pregnant sows (Schipper 1961). This study was severely limited by lack of exposure data, unequal duration of exposure between treated and untreated groups, and the lack of statistical analysis of the results. No information on adverse developmental or reproductive effects in humans following exposure to coal-tar creosote was found. An increase in relative testis weight was observed in rats administered beechwood creosote in the diet for 3 months (Miyazato et al. 1981). There were no accompanying gross or histopathological lesions of the testes in these animals, so the toxicological significance of this change is not known. Based on these limited data, it is not possible to ascertain the developmental or reproductive risk to humans of exposure to creosote.

Genotoxic Effects. The genotoxic potential of coal-tar creosote has been investigated using exclusively <u>in vitro</u> assays (Bos et al. 1984a,b, 1985, 1987; Mitchell and Tajiri 1978; Simmon and Shepherd 1978). Results of these studies are summarized in Table 2-3. The available genotoxicity data indicate that creosote is an indirect mutagen (i.e., requiring the presence of an exogenous mammalian metabolic system) and induces gene mutation in bacteria and mouse lymphoma cells. The mutagenicity of creosote observed in the conventional <u>S. typhimurium</u> assay is at least partially contributed to by the

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		, Result				
End Point	Species (Test System)	With Activation A	Without	Reference		
Prokaryotic organisms:						
Gene mutation	Salmonella typhimurium (histidine auxotrophs)	+		Simmon and Shepherd 1978 <sup>a</sup>		
	S. typhimurium	<b>+</b> <b>+</b>	- No data	Bos et al. 1985, 1987 Bos et al. 1984b		
•	S typhimurium (taped-plate assay; vapor exposure)	<b>+</b>	-	Bos et al. 1985, 1987		
	Escherichia coli WP2(TK+/-), (tryptophan auxotroph)	-	-	Simmon and Shepherd 1978 <sup>a</sup>		
Mammalian cells:						
Gene mutation	Mouse lymphoma cells	, <b>+</b> ,	<u>-</u>	Mitchell and Tajiri 1978		
Mammalian body fluids:				•		
Gene mutation	S. typhimurium (rat urine sample)	+	No data	Bos et al. 1984a		
	S. typhimurium (human urine sampla; occupational exposure)		No data	Bos et al. 1984a		
	•			• *		

as. typhimurium strains TA1537, TA98 and TA100 showed increases in frameshift mutation; strain TA1535 and E. coli straub WP2 showed no increase in base-pair substitutions. Creosote P1 is a creosote mixture American Wood Preserver's Association (AWPA) specification P-1 product used for foundation piles, freshwater piles, telephone poles, utility poles, fence posts, and other land and fresh-water uses.

<sup>+ =</sup> positive result; - = negative result

PAHs such as B[a]P and benzanthracene (Bos et al. 1984b). More recently Bos et al. (1987) identified fluoranthene as one of the major volatile components of creosote responsible for the genotoxicity observed in <u>S. typhimurium</u> strains.

In a wood-preserving factory, spot samples collected from the contaminated surfaces including that of creosote-treated wood showed mutagenic activity in <u>S. typhimurium</u> strains TA98 and TA100 (Bos et al. 1984a). Despite the contamination of the work area, no increase in mutagenic activity was detected in the urine samples of workers who inhaled creosote vapors and possibly had skin contact with the residual surface creosote. This negative mutagenic response was probably a result of a low level of creosote exposure or the fact that the components present in the workers' urine were not mutagenic (Bos et al. 1984a). In the same study, mutagenic activity was detected in the urine samples of rats injected intraperitoneally with 25 mg/kg creosote in olive oil.

Coal-tar creosote (type Pl<sup>1</sup>) was mutagenic in the mouse lymphoma assay (Mitchell and Tajiri 1978). A dose-related increase in the number of forward mutations was observed in L5178Y mouse lymphoma cells following metabolic activation. Simmon and Shepherd (1978) found that creosote (type P1) produced a mutagenic dose-response and a doubling above background mutation rate in S. typhimurium strains TA1537, TA98 and TA100. S. typhimurium strain TA1535 and Escherichia coli WP2 strain did not demonstrate a positive response with metabolic activation. These results indicate that the genetic mode of action of creosote in S. typhimurium is by frameshift mutation. In recent studies Bos et al. (1984b) examined the mutagenic potential of several fractions of creosote (type P1) in strains TA98 and TA1537. The mutagenicity of creosote was found to be associated at least partially with PAHs (B[a]P and benzanthracene) which were detected in concentrations of 0.18% and 1.1%, respectively (Bos et al. 1983, 1984b). Both compounds, although mutagenic in conventional Ames assay, showed no activity in a S. typhimurium taped-plate assay, probably because of their low volatility (Bos et al. 1985). When creosote samples were further tested for the presence of "volatile mutagens," vapors escaping from creosote increased the number of revertants per plate in the presence of exogenous mammalian metabolic activation system (S9 mix) (Bos et al. 1985). Results of the study suggest that the volatile components of creosote may also contribute to the genotoxic risk from occupational exposure to creosote. In a recent study, Bos et al. (1987) found that creosote contained mutagens which were volatile at 37.C. These were present in the

<sup>&</sup>lt;sup>1</sup>Creosote is classified according to a set of standards regarding physical property specifications for creosote that must be met for certain uses of creosote. These specifications are presented in Table 3-4. Type P1/P13 creosote is straight creosote distillate and is used in ground contact, land, and fresh and marine water applications. Type P2 creosote is used by the railroad industry in the treatment of railroad crossties.

distillation fraction having the highest boiling range (more than 360·C). Upon further high-pressure liquid chromatography (HPLC) and ultraviolet spectroscopic analysis of the fraction and mutagenicity tests, the mutagenic response observed was correlated with the presence of fluoranthene (5.2%) in the creosote samples. The commercially available fluoranthene also tested positive in the taped-plate assay. The available genotoxicity data suggest that creosote has a potential to induce gene mutation in humans exposed via inhalation, and perhaps other routes as well, at hazardous waste sites. However, much of the data discussed above describes the mutagenic activity of selected PAHs. It is not yet known how the various PAHs contained in the coal-tar creosote mixture interact, and what the effects of these possible interactions are on the ultimate genotoxic expression of the mixture.

Cancer. Various case reports and the results of cross-sectional occupational surveys associate chronic occupational creosote exposure with the development of skin cancer (Cookson 1924; Henry 1946, 1947; Lenson 1956; O'Donovan 1920). These papers reported a similar disease etiology for different groups of workers exposed to creosote that included the development of dermatoses, such as squamous papillomas, that progressed to carcinoma, usually squamous-cell carcinoma. The latency period for the development of dermatoses, such as squamous papillomas, was usually 20-25 years (Cookson 1924; Henry 1946, 1947; O'Donovan 1920). The latency period for the development of carcinoma was on average 25 to 30 years (Cookson 1924; Henry 1946, 1947; O'Donovan 1920). Worker exposure in the past was much greater than it now is because of less-sophisticated industrial practices used in the past, the lack of knowledge concerning occupational hygiene, and the current recognition of the dangers of excessive exposure to the health of workers. Other factors that should be considered when extrapolating the findings of this older literature to present conditions are the role of exposure to ultraviolet radiation in the form of sunshine in these workers (ultraviolet radiation is now known to be a major cause of skin cancer), the composition of the creosote products, and other health factors that differ in Great Britain prior to 1940 and the present. Although these studies lack information concerning specific exposures and did not consider other risk factors for the development of skin cancer, when taken as a group, and in view of the presence of carcinogenic PAHs, they present convincing evidence of a relationship between chronic creosote exposure and the development of skin carcinoma in humans.

On the other hand, several studies are available that suggest that there is no association between exposure to creosote or other coal-tar products and cancer in humans. No increased incidence of non-melanoma skin cancer was observed in 426 patients 25 years after they had undergone 4 years of coal-tar medicinal therapy in combination with ultraviolet light for the treatment of atopic dermatitis and neurodermatitis (Maugham et al. 1980). This study is limited in that follow-up occurred in only 72% of the patients, and there was no discussion of recall bias or of the effects of mobility (i.e. relocation of

the study participants) on the results. In a 25-year retrospective study of 280 psoriatic patients treated with crude coal-tar in combination with ultraviolet radiation at the Mayo clinic, it was found that the incidence of skin cancer in these patients was not significantly increased above the expected incidence for the general population (Pittelkow et al. 1981). In investigations on populations using therapeutic coal-tar skin products, exposure was to a different coal tar mixture. Creosote is a lower boiling distillate fraction of crude coal-tar than the components contained in the skin products, and does not contain the concentrations of putative carcinogenic PAH compounds that are present in the tars (Wright et al. 1985).

Another study was located that reported that there was no increase in the risk of skin, bladder, or lung cancer in wood treatment plant workers (Tabershaw Occupational Medicine Associates 1980). Limitations in this study include: the study population was small; the study population was comprised of 46.5% blacks who experience a very low incidence of skin cancer as compared to whites, thus biasing the results; the exposure and follow-up periods did not allow a long enough latency period for tumor development; and there was no verification provided that those studied were actually exposed to creosote or coal-tar. However, it is possible that because black people are less sensitive to ultraviolet light-induced cancer, their prominence in this study might serve to help isolate the effects of creosote alone from those of ultraviolet light or ultraviolet light plus creosote.

A large body of evidence exists indicating that coal-tar creosote and several of its fractions are carcinogenic when applied to the skin of laboratory animals. Many of the early studies that reported that coal-tar creosote is carcinogenic when dermally applied to rodent skin are limited in that they lack appropriate negative control data, the dose of creosote and the chemical composition of the fractions studied were not quantified, and no other tissues were examined. The results from later studies that include appropriate control groups are consistent with the findings of the earlier studies. Dermally applied creosote can also act as a tumor-initiating agent when applied prior to croton oil treatment, and can enhance and accelerate tumor induction by B[a]P. IARC (1985) has classified creosote as a Group 2A, probable human carcinogen based on limited human evidence and sufficient animal evidence of carcinogenicity. Thus, based on the human and animal data summarized above, it is possible that individuals who chronically come in skin contact with creosote may be at an elevated risk for developing skin cancer, particularly when exposure to other carcinogenic substances occurs, as is a likely scenario in areas surrounding hazardous waste sites.

# 2.5 BIOMARKERS OF EXPOSURE AND EFFECT

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Biomarkers are broadly defined as indicators signaling events in biologic systems or samples. They have been classified as markers of exposure, markers of effect, and markers of susceptibility (NAS/NRC 1989).

A biomarker of exposure is a xenobiotic substance or its metabolite(s) or the product of an interaction between a xenobiotic agent and some target molecule or cell that is measured within a compartment of an organism (NAS/NRC 1989). The preferred biomarkers of exposure are generally the substance itself or substance-specific metabolites in readily obtainable body fluid or excreta. However, several factors can confound the use and interpretation of biomarkers of exposure. The body burden of a substance may be the result of exposures from more than one source. The substance being measured may be a metabolite of another xenobiotic substance (e.g., high urinary levels of phenol can result from exposure to several different aromatic compounds). Depending on the properties of the substance (e.g., biologic half-life) and environmental conditions (e.g., duration and route of exposure), the substance and all of its metabolites may have left the body by the time biologic samples can be taken. It may be difficult to identify individuals exposed to hazardous substances that are commonly found in body tissues and fluids (e.g., essential mineral nutrients such as copper, zinc, and selenium). Biomarkers of exposure to creosote are discussed in Section 2.5.1.

Biomarkers of effect are defined as any measurable biochemical, physiologic, or other alteration within an organism that, depending on magnitude, can be recognized as an established or potential health impairment or disease (NAS/NRC 1989). This definition encompasses biochemical or cellular signals of tissue dysfunction (e.g., increased liver enzyme activity or pathologic changes in female genital epithelial cells), as well as physiologic signs of dysfunction such as increased blood pressure or decreased lung capacity. Note that these markers are often not substance specific. They also may not be directly adverse, but can indicate potential health impairment (e.g., DNA adducts). Biomarkers of effects caused by creosote are discussed in Section 2.5.2.

A biomarker of susceptibility is an indicator of an inherent or acquired limitation of an organism's ability to respond to the challenge of exposure to a specific xenobiotic substance. It can be an intrinsic genetic or other characteristic or a preexisting disease that results in an increase in absorbed dose, biologically effective dose, or target tissue response. If biomarkers of susceptibility exist, they are discussed in Section 2.7, "POPULATIONS THAT ARE UNUSUALLY SUSCEPTIBLE."

# 2.5.1 Biomarkers Used to Identify or Quantify Exposure to Creosote

No method is currently available to measure the parent crossote mixture in human tissues or fluids. However, the PAH components of the crossote mixture and their metabolites can be measured in the urine of exposed individuals. For example, Jongeneelen et al. (1985) found a metabolite of crossote, 1-hydroxypyrene, in concentrations of 1-40  $\mu$ g/g creatinine in urine samples taken from workers who handled approximately 2,400 g crossote/day. The amount of 1-hydroxypyrene detected in urine samples taken during the weekend was less than that detected during the weekdays, when the exposure was

presumably higher than on the weekends. No correlation was found between occupational exposure levels and urine levels, so it is not known whether urine metabolites could be detected following exposure to low levels of creosote. The identification of PAH metabolites in the urine could serve as a method of biological monitoring of exposed workers, and possibly individuals living in the vicinity of hazardous waste sites where creosote has been detected following both short- and long-term exposure. However, because of the ubiquitous nature of PAHs in the environment, detection of PAH metabolites in the body tissues or fluids is not specific for exposure to creosote. PAH exposure can occur from a variety of sources, and there is no way to determine if creosote was the source.

PAHs form DNA adducts that can be measured in body tissues or blood following exposure to creosote that contains PAHs. Again, these PAH-DNA adducts are not specific for coal-tar creosote, and the adducts measured could have been from exposure to other sources of PAHs.

# 2.5.2 Biomarkers Used to Characterize Effects Caused by Creosote

The available genotoxicity data indicate that creosote is an indirect mutagen (i.e., requiring the presence of an exogenous mammalian metabolic system) and induces gene mutation in bacteria and mouse lymphoma cells. The mutagenicity of creosote observed in the conventional S. typhimurium assay is at least partially contributed to by the PAHs such as B[a]P and benzanthracene. However, because these results are exclusively from in vitro tests and the limited genotoxicity tests conducted on urine obtained from humans exposed to creosote have been negative, these changes cannot be considered specific biomarkers of effects caused by creosote, nor is it possible to distinguish whether the genotoxic effects result from either acute or chronic exposure to either low or high levels of creosote because all of the data were from in vitro studies. Furthermore, because the mutagenicity of creosote is at least partially due to its PAH components, exposure to PAHs from other sources could produce the same results. Coal-tar creosote exerts its acute toxic effects primarily via dermal exposure, causing architectural damage to the tissues with which it comes in contact. Therefore, burns and irritation of the skin and eyes are the most frequent manifestations of coaltar creosote toxicity following acute dermal exposure to high levels. However, damage to the skin is not specific to creosote, and can be seen with other corrosive or photosensitizing agents. No other biomarkers (specific or otherwise) have been identified following exposure to creosote.

## 2.6 INTERACTIONS WITH OTHER CHEMICALS

The primary interactions known to occur between coal-tar creosote and other substances involve the induction of cancer. Coal-tar creosote is a complex mixture of organic substances consisting predominantly of liquid and solid aromatic hydrocarbons. Several of these components of coal-tar creosote

are known carcinogens as well as cocarcinogens, initiators, promoters, potentiators, or inhibitors of carcinogenesis. As discussed in Section 2.2.3.8 (Dermal Exposure, Cancer), coal-tar creosote and several of its fractions are carcinogenic when applied to the skin of mice. Dermally applied creosote can also act as a tumor-initiating agent when applied prior to croton oil treatment, and can enhance and accelerate tumor induction by B[a]P. Thus, the risk of cancer following dermal exposure to creosote is likely to be enhanced when concurrent exposure to other potential cocarcinogens, tumor promoters, initiators, and potentiators occurs. Due to the ubiquitous nature of PAHs and other carcinogenic substances in the environment, particularly at hazardous waste sites, the likelihood that these types of synergistic interactions with creosote will occur is good.

Pentachlorophenol and arsenical compounds are also used in wood preserving. For this reason, it is likely that they will be found with creosote at hazardous waste sites. However, there is no information available on the potential interactions of creosote with pentachlorophenol or arsenical compounds.

## 2.7 POPULATIONS THAT ARE UNUSUALLY SUSCEPTIBLE

Data indicate that the population in general may be at increased risk of developing skin cancer following prolonged dermal exposure to creosote. Data also indicate that subsections of the human population may be unusually susceptible to the toxic effects of creosote. These include people with a history of excessive sun exposure, people with skin diseases, and people with exposure to other substances that act as cocarcinogens, tumor promoters, initiators, and potentiators. There is some limited evidence, based on animal studies and the known health effects of the PAH constituents of creosote, that indicates that additional subsections of the population may be susceptible to the toxic effects of creosote. These include fetuses, people with kidney or liver disease, people with deficient immune systems due to disease or advanced age, and people with the genetic trait of inducible aryl hydrocarbon hydroxylase (AHH). AHH is an enzyme contained in the microsomal fraction of the cell that is responsible for the biotransformation of aryl compounds, such as PAHs. When this enzyme is induced, the rate at which aryl compounds are biotransformed into toxic intermediates is increased, rendering these individuals at higher risk. Genetic polymorphism does exist in the general population with regard to the rates of aryl hydrocarbon hydroxylation (i.e., differences in the P-450 and P-448 enzymes).

The results of occupational studies (Henry 1946, 1947), case reports (Cookson 1924; Lenson 1956; O'Donovan 1920), and experimental animal studies (Boutwell and Bosch 1958; Poel and Kammer 1957; Roe et al. 1958) indicate that the general population may be at risk of developing skin cancer following prolonged dermal exposure to creosote. People with skin damaged from excessive sun exposure, disease, or exposure to other substances that potentiate the carcinogenic effect of creosote (Cabot et al. 1984; Lenson

1956; Lijinsky et al. 1956; Sall and Shear 1940) are at an increased risk of developing skin disease, including cancer, from creosote exposure.

Limited evidence from experimental studies of pregnant pigs (Schipper 1961) and pregnant rats (Srivastava et al. 1986) indicate that components of creosote, such as PAHs, cross the placenta following maternal inhalation and may cause fetotoxic effects. Data from a human case report (Bowman et al. 1984) and experimental animal studies (Miyazato et al. 1981, 1984a,b) indicate that kidney or liver disease may increase the risk of developing adverse health effects following ingestion of creosote.

Two other subpopulations may be susceptible to certain components of creosote, such as PAHs; one would be people with deficient immune systems as a result of disease or advanced age, and the second would be people who carry the genetic trait of inducible AHH, one of the mixed-function oxidases. Calabrese (1978) reported that people with deficient immune systems may be at high risk of developing adverse health effects due to exposure to carcinogens, such as PAHs (Stjernsward 1966, 1969; Szakal and Hanna 1972). PAHs are metabolized to reactive intermediates, which are thought to cause cell transformation, mutagenicity, and cytotoxicity. It has been proposed that genetically expressed AHH inducibility is related to the development of bronchogenic carcinoma in persons exposed to PAHs contained in tobacco smoke. Approximately 45% of the general population are considered to be at high risk, and 9% of the 45% are considered to be at very high risk of developing bronchogenic carcinoma following exposure to PAHs (Calabrese 1978). These percentages were estimated from the population frequency of genetically controlled AHH induction (Calabrese 1978).

## 2.8 ADEQUACY OF THE DATABASE

Section 104(i)5 of CERCIA directs the Administrator of ATSDR (in consultation with the Administrator of EPA and agencies and programs of the Public Health Service) to assess whether adequate information on the health effects of creosote is available. Where adequate information is not available, ATSDR, in cooperation with the National Toxicology Program (NTP), is required to assure the initiation of a program of research designed to determine the health effects (and techniques for developing methods to determine such health effects) of creosote.

The following categories of possible data needs have been identified by a joint team of scientists from ATSDR, NTP, and EPA. They are defined as substance-specific informational needs that, if met would reduce or eliminate the uncertainties of human health assessment. In the future, the identified data needs will be evaluated and prioritized, and a substance-specific research agenda will be proposed.

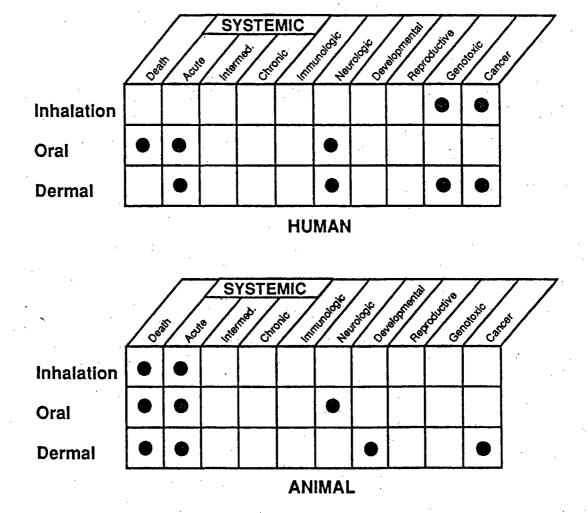
# 2.8.1 Existing Information on the Health Effects of Creosote

The existing data on health effects of inhalation, oral, and dermal exposure of humans and animals to creosote are summarized in Figure 2-3. The purpose of this figure is to illustrate the existing information concerning the health effects of creosote. Each dot in the figure indicates that one or more studies provide information associated with that particular effect. The dot does not imply anything about the quality of the study or studies. Gaps in this figure should not be interpreted as "data needs" information. The vast majority of literature reviewed concerning the health effects of coal-tar creosote in humans described case reports of workers exposed predominantly via the dermal route, with a few older reports describing the consequences of accidental or intentional ingestion of creosote. Despite the limitations inherent in these types of studies, an association between dermal exposure to coal-tar creosote and dermal irritation, phototoxicity, and cancer in humans is evident.

The database for the health effects of creosote in experimental animals is lacking, and consists primarily of acute lethality studies or old animal studies that would be considered inadequate by current standards. The systemic effects of ingested creosote have only been well described for beechwood creosote. Little information is available on the effects of creosote following inhalation exposure. However, dermal exposure to coal-tar creosote has been shown in numerous studies to induce skin and lung tumors in mice. Since coal-tar creosote is a complex mixture consisting primarily of PAHs, the toxic effects of coal-tar creosote may be inferred from available information on these constituents. However, given the fact that many of these constituents are known cocarcinogens, initiators, promoters, and potentiators of carcinogenesis, the possibility for the occurrence of synergistic interactions in creosote cannot be ruled out. Thus, information on the toxicity of the various components of coal-tar creosote cannot take the place of sound data on the toxic effects of the creosote mixture itself.

### 2.8.2 Identification of Data Needs

Acute-Duration Exposure. Information is available on the effects of acute-duration exposures to creosote in humans and animals (oral and dermal). The type of information available includes  ${\rm ID}_{50}{\rm S}$  and acute toxicity in animals (coal-tar and beechwood creosote), and acute toxicity following accidental or intentional ingestion or dermal exposure in humans (coal-tar creosote). Coaltar creosote exerts its acute toxic effects primarily via dermal exposure, causing architectural damage to the tissues with which it it comes in contact, such as the skin and eyes. Acute ingestion of coal-tar creosote appears to affect primarily the kidney and liver. Thus, the toxic effects of coal-tar creosote on the skin following single-dose dermal exposure in humans are well characterized, but little else is known regarding the systemic effects of this form of creosote in either humans or animals. The available information is



# Existing Studies

FIGURE 2-3. Existing Information on Health Effects of Coal Tar Creosote

insufficient to derive either an acute oral or inhalation MRL for coal-tar creosote because human reports that identified target organs lacked exposure information, and no short-term animal studies exist that describe effects other than death. Identification of target organs from short-term animal studies following oral and dermal exposure would be useful in assessing the risk associated with the acute ingestion or skin contact by humans of coal-tar creosote contaminated water or soils. The pharmacokinetic data on coal-tar creosote are insufficient to determine whether similar effects may be expected to occur across different routes of exposure. However, since creosote appears to cause route-of-entry adverse effects (e.g., damage to the skin following dermal contact), it may not be possible to predict effects following exposure by one route based on effects observed following exposure by another route.

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Intermediate-Duration Exposure. Information is available on the effects of intermediate-duration dermal exposures to coal-tar creosote in humans and the effects of intermediate-duration exposures to beechwood creosote (oral) and coal-tar creosote (dermal) in animals. The exact duration and level of exposure in the human studies generally cannot be quantified because the information is derived from anecdotal case reports rather than controlled epidemiological studies. The animal studies with beechwood creosote describe predominantly hepatic and renal end points, and those conducted with coal-tar creosote describe dermal, but very rarely systemic, effects. Little or no reliable information on respiratory, cardiovascular, gastrointestinal, hematological, or musculoskeletal effects in animals is available. available information is insufficient to derive either an intermediate oral or inhalation MRL for coal-tar creosote because no intermediate-duration animal studies exist that describe effects other than on the skin. Given the widespread use of coal-tar creosote as a wood preservative, and the fact that beechwood creosote is rarely used today, more information on the systemic effects of intermediate-duration exposures to coal-tar creosote by the oral and dermal routes (by conducting 90-day subchronic toxicity studies) would be useful to identify target organs in animals in order to assess the risk associated with the intermediate-duration ingestion or skin contact by humans of coal-tar creosote contaminated water or soils. The pharmacokinetic data on coal-tar creosote are insufficient to determine whether similar effects may be expected to occur across different routes of exposure. However, since creosote appears to cause route-of-entry adverse effects (e.g., damage to the skin following dermal contact), it may not be possible to predict effects following exposure by one route based on effects observed following exposure by another route.

Chronic-Duration Exposure and Cancer. Information is available on the effects of chronic-duration dermal exposures to coal-tar creosote in humans and the effects of chronic-duration exposures to beechwood creosote (oral) and coal-tar creosote (dermal) in animals. The exact duration and level of exposure in the human studies generally cannot be quantified because the information is derived from anecdotal case reports rather than controlled

epidemiological studies. The animal studies with beechwood creosote describe predominantly hepatic and renal end points, and those conducted with coal-tar ereosote describe dermal, but very rarely systemic, effects. Little or no reliable information on respiratory, cardiovascular, gastrointestinal, hematological, or musculoskeletal effects in animals is available. available information is insufficient to derive either a chronic oral or inhalation MRL for coal-tar creosote because no chronic-duration animal studies exist that describe effects other than on the skin. Given the widespread use of coal-tar creosote as a wood preservative, and the fact that beechwood creosote is rarely used today, more information on the systemic effects of chronic-duration exposures to coal-tar creosote by the oral and dermal routes would be useful to identify target organs in animals in order to assess the risk associated with the chronic-duration ingestion or skin contact by humans of coal-tar creosote contaminated water or soils. pharmacokinetic data on coal-tar creosote are insufficient to determine whether similar effects may be expected to occur across different routes of exposure. However, since creosote appears to cause route-of-entry adverse effects (e.g., damage to the skin following dermal contact), it may not be possible to predict effects following exposure by one route based on effects observed following exposure by another route.

Various older case reports and the results of cross-sectional occupational surveys associate chronic occupational creosote exposure with the development of skin cancer. However, several newer studies are available that suggest that there is no association between exposure to creosote or other coal-tar products and cancer in humans. Several skin painting studies have been conducted in animals using coal-tar creosote and its various fractions. Although many of these studies would be considered inadequate by current standards, the results nevertheless indicate that coal-tar creosote and its constituents can induce skin tumors as well as act as tumor initiators and promoters. Although carcinogenicity studies have been conducted with beechwood creosote by the oral route, the relevance of the findings to coaltar creosote is not known. More information on the carcinogenic potential of chronically ingested coal-tar creosote (e.g., an oral bioassay) would be useful. The pharmacokinetic data on coal-tar creosote are insufficient to determine whether similar effects may be expected to occur across different routes of exposure. However, since creosote appears to cause route-of-entry adverse effects (e.g., skin tumors following dermal contact), it may not be possible to predict effects following exposure by one route based on effects observed following exposure by another route.

Genotoxicity. The genotoxic potential of coal-tar creosote has been investigated almost exclusively using <u>in vitro</u> assays. The limited genotoxicity tests that have been conducted on urine obtained from humans exposed to creosote have been negative. The available data indicate that creosote is an indirect mutagen and induces gene mutation in bacteria and mouse lymphoma cells. However, a substantial database exists on the genotoxic

effects of the PAHs found in the creosote mixture. More <u>in vivo</u> assays with creosote would be useful to more completely characterize the genotoxic potential of the creosote mixture.

Reproductive Toxicity. No information on the reproductive effects of coal-tar creosote in humans or animals is available. However, limited data available on benzo(a)pyrene indicate that this PAH has a potential to induce adverse reproductive effects. An increase in relative testis weight was observed in rats administered beechwood creosote in the diet for 3 months. There were no accompanying gross or histopathological lesions of the testes in these animals, so the toxicological significance of this change is not known. Given the widespread potential for exposure to coal-tar creosote, and the indication that certain PAH components of coal-tar creosote as well as beechwood creosote may be reproductive toxicants, reproductive organ pathology she id be examined in any 90-day studies that may be conducted. If these winary data indicate that creosote may present a reproductive hazard, to a multi-generation reproductive toxicity studies should be conducted by the onel and dermal routes of exposure. The pharmacokinetic data on coal-tar cheosope are insufficient to determine whether similar effects may be expected the accour across different routes of exposure. However, since creosote appears to cause route-of-entry adverse effects (e.g., skin tumors following dermal manufacture, it may not be possible to predict effects following exposure by one haned on effects observed following exposure by another route.

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Developmental Toxicity. Information on the developmental effects of eccesote in humans was not found. Data from one severely limited study in pigs indicate that coal-tar creosote has the potential to be fetotoxic. Limited information available on a few PAHs (mostly benzo[a]pyrene) indicate that information available on a few PAHs (mostly benzo[a]pyrene) indicate ingested PAHs have a potential to induce adverse developmental effects in animals. Given the widespread potential for exposure to coal-tar creosote, and the suggestive data mentioned above, oral and dermal developmental toxicity studies in animals would be useful to assess the potential risk for sote-induced adverse developmental effects. The pharmacokinetic data on coal-tar creosote are insufficient to determine whether similar effects may be expected to occur across different routes of exposure. However, since creosote appears to cause route-of-entry adverse effects (e.g., skin tumors owing dermal contact), it may not be possible to predict effects following sure by one route based on effects observed following exposure by another route.

Immunotoxicity. The only available information on the immunological effects of creosote describes the occurrence of acute allergic dermatitis following exposure to creosote bush resin. The relevance of these findings to exposure to coal-tar creosote is not known. However, since creosote bush resin and coal-tar creosote do contain some similar constituents (e.g., phenolic derivatives), and people with deficient immune systems may be at high risk of developing adverse health effects due to exposure to carcinogens, such

as the PAHs contained in creosote, more information on the immunologic effects of dermally applied coal-tar creosote would be useful to better characterize the toxic effects of exposure to coal-tar creosote. Preliminary information on the potential for creosote to induce immunotoxic effects may be obtained from 90-day studies that examine effects on lymphoid tissue and blood components. If adverse effects on these parameters are observed, then a full battery of immunotoxic tests may be warranted to further characterize the potential immunotoxicity of creosote.

Neurotoxicity. The available information describes neurological involvement in humans and animals following short-term high-level oral exposure to creosote (both beechwood and coal-tar). These effects are generally excitatory in nature (i.e., convulsions). No information is available on the long-term neurotoxic effects of low-level exposure to coal-tar creosote in humans and animals. Long-term neurotoxicity studies in animals, using sensitive functional and neuropathological tests, would be useful in determining if coal-tar creosote is a neurotoxic agent.

Epidemiological and Human Dosimetry Studies. Few controlled epidemiological studies have been conducted in humans on the effects of exposure to coal-tar creosote. In particular, epidemiological studies of workers in creosote treatment plants would be useful to more fully assess the risk of inhalation and dermal exposure to coal-tar creosote. Most of the available information on the effects of coal-tar creosote in humans comes from cases of acute poisoning following the accidental or intentional exposure to coal-tar creosote and from occupational exposures in the wood-preserving and construction industries. Limitations inherent in these studies include unquantified exposure concentrations and durations, as well as concomitant exposure to other potentially toxic substances. The few available industrial surveys and epidemiological studies are limited in their usefulness because of small sample size, short follow-up periods, and brief exposure periods. Despite their inadequacies, studies in humans suggest that coal-tar creosote is a dermal irritant and a carcinogen following dermal exposure. Wellcontrolled epidemiological studies of people living in close proximity to areas where coal-tar creosote has been detected in surface and groundwater, near hazardous waste sites, and of people occupationally exposed could add to and clarify the existing database on creosote-induced human health effects. Particular effects to be examined include cancer (of the skin and other organs) and other adverse skin effects.

Biomarkers of Exposure and Effect. No method is currently available to measure the parent creosote mixture in human tissues or fluids. However, the PAH components of the creosote mixture and their metabolites (e.g., 1-hydroxypyrene) can be measured in the urine of exposed individuals following relatively high-level exposures of acute and chronic duration. The identification of PAH metabolites in the urine could serve as a method of biological monitoring of exposed workers, and possibly individuals living in

the vicinity of hazardous waste sites where creosote has been detected. However, because of the ubiquitous nature of PAHs in the environment, detection of PAH metabolites in the body tissues or fluids is not specific for exposure to creosote. PAH exposure can occur from a variety of sources, and there is no way to determine if creosote was the source. PAHs form DNA adducts that can be measured in body tissues or blood following exposure to creosote containing PAHs. Again, these PAH-DNA adducts are not specific for coal-tar creosote, and the adducts measured could have been from exposure to other sources of PAHs. Therefore, a biomarker of exposure specific to creosote would be useful to monitor exposure to this mixture.

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The formation of benzo(a)pyrene-DNA adducts has been demonstrated and may also serve as a biomarker of PAH-induced carcinogenicity. However, these adducts are not specific for coal-tar creosote exposure, as exposure to benzo(a)pyrene from sources other than coal-tar creosote can occur. Studies to identify and measure effects more diagnostic of coal-tar creosote-specific injury would be useful. Also, increasing the sensitivity of these tests would be valuable in evaluating the health status of individuals who have been exposed to low levels of creosote.

Absorption, Distribution, Metabolism, and Excretion. Studies monitoring the pharmacokinetics of the coal-tar creosote mixture are limited. Much of the information regarding the disposition of creosote is based on indirect evidence or the pharmacokinetic information available on a single class of components of creosote, the PAHs.

Absorption of creosote occurs following all routes of exposure. The presence of creosote components in tissues and the presence of metabolites in urine were evidence of its absorption. However, no studies are available that quantify the extent and rate of creosote absorption.

Studies in humans or animals regarding the distribution of creosote are not available. Its distribution is based on assumptions derived from studies that monitored the distribution of PAHs, components of creosote.

The metabolism of creosote has not been extensively studied, but preliminary results indicate that hydroxylation of the major PAH components is a principal degradation pathway in both humans and animals following all routes of exposure. 1-Hydroxypyrene is one metabolite that has been identified, but there were no studies available regarding the identification of other metabolites. Elucidation of additional biotransformation pathways and products is also important in examining potential toxic effects of creosote. Moreover, no studies were located regarding the rate or extent of creosote metabolism.

Studies regarding the excretion of creosote by humans or animals were not available. It is known that PAHs and their metabolites are primarily excreted

in the bile and the feces. However, direct excretion studies with creosote would be more useful. Information is available regarding the disposition of creosote's individual components, but no information is available regarding how these components interact to affect the overall disposition.

In summary, no data are available regarding the toxicokinetics of the creosote mixture and all information must currently be inferred from what is known about the PAH components of creosote. Interactions between the components of the creosote mixture could occur that could alter the rate and extent of absorption, distribution, metabolism, and excretion of creosote from what might be predicted based on what is known about the individual PAH components. Therefore, more information on the toxicokinetics of the creosote mixture itself would be useful to predict possible target organs of toxicity as well as allow for extrapolation of toxic effects across routes of exposure.

Comparative Toxicokinetics. The available information indicates that the absorption, distribution, metabolism, and excretion of creosote is qualitatively similar in humans and rodents. This general conclusion was primarily based on evidence derived from studies on the individual PAH components of creosote. Detailed pharmacokinetic studies in humans and animals specific to the creosote mixture would provide a better indication of species differences and indicate whether the ability to extrapolate across species may be possible in the future.

# 2.8.3 On-going Studies

Creosote is currently subject to an EPA Federal Insecticide, Fungicide, and Rodenticide Act (FIFRA) registration standard and data call-in, and the Creosote Council II is planning to conduct a research program that includes testing in subchronic inhalation, subchronic dermal, developmental, and reproductive toxicity.

#### 3.1 CHEMICAL IDENTITY

The chemical synonyms and identification numbers for coal-tar creosote are listed in Table 3-1. Creosote is a complex mixture of variable composition containing primarily aromatic ring compounds and condensed aromatic ring compounds. Therefore, it is not possible to represent the chemical formula and structure of the creosote mixture.

# 3.2 PHYSICAL AND CHEMICAL PROPERTIES

Important physical and chemical properties of coal-tar creosote are listed in Table 3-2.

Coal-tar creosote is defined by the American Wood Preserver's Association (1988) as:

A distillate derived from coal-tar. As used in the wood preserving industry, creosote denotes a distillate of coal-tar produced by the high temperature carbonization of bituminous coal. Creosote consists principally of liquid and solid aromatic hydrocarbons and contains some tar acids and tar bases; it is heavier than water and has a continuous boiling range beginning at about 200°C.

It is thus a complex mixture typically composed of 85% PAHs and 2%-17% phenolics (Bedient et al. 1984). The composition of the mixture may also vary across lots and across manufacturers. The major PAH components of creosote are listed in Table 3-3. Creosote is further classified according to a set of standards regarding physical property specifications for creosote that must be met for certain types of creosote. The physical property specifications for the different types of creosote are listed in Table 3-4. Type P1/P13 creosote is a straight creosote distillate and is used in ground contact, land, and fresh and marine water applications. Type P2 creosote is used by the railroad industry in the treatment of railroad crossties.

Coal-tar itself is produced by the carbonization, or coking of coal. Coal-tar is defined by Hawley (1977) as:

A black, viscous liquid (or semi-solid), naphthalene-like odor, sharp burning taste; obtained by the destructive distillation of bituminous coal, as in coke ovens; 1 ton of coal yields 8.8 gallons of coal-tar. Combustible. Specific gravity 1.18-1.23 (66/60°F). Soluble in ether, benzene, carbon disulfide, chloroform; partially soluble in alcohol, acetone, methanol, and benzene; only slightly soluble in water.

To distinguish another product of coal-tar from creosote, coal-tar neutral oils are usually defined as a mixture of naphthalene, fluorene, anthracene, and other neutral hydrocarbons (Smale, 1977). Neutral hydrocarbons consist of those coal-tar hydrocarbons other than coal-tar acids (such as phenols, cresols, and cresylic acids), and coal-tar bases (such as pyridines, quinolines, and acridines).

Because of the variability in feedstock and manufacturing processes, presentation of exact values for various properties presented in Table 3-2 are not possible.

TABLE 3-1. Chemical Identity of Coal-tar Creosote.

	Value	Reference
Chemical name	Coal-tar creosote	AWPA 1988
Synonyms	Creosote Oil, Dead Oil, Brick Oil, Coal-tar Oil, Creosote Pl, Heavy Oil, Liquid Pitch Oil, Wash Oil	Windholz 1983; Weiss 1986; HSDB 1988
Trade Names	Preserv-O-Sote	HSDB 1988
Chemical Formula	No data <sup>a</sup>	Windholz 1983 HSDB 1988
Chemical Structure	No data <sup>e</sup>	
Wiswesser Line Notation	UVCB	HSDB 1988
Identification Numbers	<b>5:</b>	
CAS Registry NIOSH RTECS EPA Hazardous Waste OHM/TADS DOT/UN/NA/IMCO shipping HSDB	8001-58-9 GF8615000 U051 No data DOT: 1933; IMCO: 3.2; IMCO/UN: 9/1993 6299	Weiss 1986 HSDB 1988 HSDB 1988 Weiss 1986 HSDB 1988
NCI	No data	•

<sup>&</sup>lt;sup>a</sup> Creosote is a mixed compound composed primarily of polycyclic aromatic hydrocarbons including phenanthrene, acenaphthene, fluorene, anthracene, and pyridine.

AWPA = American Wood Preservers' Association

CAS = Chemical Abstracts Service

NIOSH = National Institute for Occupational Safety and Health

RTECS = Registry of Toxic Effects of Chemical Substances

DOT/UN/NA/IMCO = Department of Transportation/United Nations/North

America/International Maritime Dangerous Goods Code

HSDB = Hazardous Substances Data Bank

TABLE 3-2. Physical and Chemical Properties of Coal-Tar Creosote

Property	Value	Reference
Molecular weight	No data	
Color	Translucent amber to	Windholz 1983
60101	brown to black	WINDING 1903
	liquid	
Physical state	Liquid	Weiss 1988
Melting point	No data	#6135 1700
Boiling point	>180° C	Weiss 1986
borring points	194-400° C	HSDB 1988
Density (g/cm3)	1.07 (20° C)	DOT 1985
Odor	Sharp odor	Windholz 1983
	Tarry, aromatic odor	Weiss 1986
•	Smoky odor	HSDB 1988
Odor threshold:	•	,
Water	No data	
Air	No data	*
Solubility:		
Water	Practically insoluble	Windholz 1983
Òrganic solvents	Miscible with alcohol,	HSDB 1988
	ether, fixed or volatile	•
•	oils	
Partition coefficients	No data	
Vapor pressure	No data	
Henry's law constant	No data .	•
Autoignition Temperature	335° C	Windholz 1983
Flashpoint	75° C	Windholz 1983
Flammability limits	1.3-8%	HSDB 1988

TABLE 3-3. Major Chemical Components of Typical Wood Preservative Creosote\*

	Percent			
Component P	1/P13 Creosote*	* P2 Creosote		
Indene	3.3	0.5		
Methylbenzofurans and				
xylenol	1.0			
Naphthalene	14.3	8.7		
Quinoline	0.9	0.7		
2-Methylnaphthalene ar	nd	•		
indole	4.9	5.6		
1-Methylnaphthalene	2.1	2.1		
Biphenyl	1.3	1.5		
Dimethylnaphthalenes	2.1			
Acenapthylene	0.8			
Acenaphthene	4.1	4.9		
Dibenzofuran	3.5	3.4		
Fluorene	4.7	4.5		
Methyldibenzofurans	3.0			
Methylfluorenes	2.2	0.8		
Dibenzothiophene	0.9			
Phenanthrene	12.7	16.9		
Anthracene	5.6	5.3		
Benzoquinolines	0.8			
Carbazole	0.9	2.2		
2-& 3-Methylphenanthre	enes 2.0	3.0		
Fluoranthene	6.0	7.3		
Pyrene	5.0	6.5		
Benzofluorenes and	•			
methylfluoranthenes	3.0			
Methylpyrenes	1.9			
Chrysene	1.5	3.5		
Methylchrysene and				
benzanthracenes	1.0	1.1		

<sup>\*</sup>These values are calculated averages which are representative of creosote. Actual concentrations of specific constitutents can vary. Naphthalene, for example, can range from five to twenty percent.

Source: Butala 1990

<sup>\*\*</sup>See Table 3-4 for the physical property specifications for each type of creosote.

TABLE 3-4. Physical Property Specifications for Various Types of Creosote

	Range of Properties			
Typical Physical Properties	Creosote Distillate (Pl/Pl3)	Creosote/ Coal-tar Solution (P2)		
Water, Percent by Volume	0.0-0.7		0.2-0.13	
	·			
Specific Gravity at 38°C:			•	
Whole Fraction	1.059-1.102		1.083-1.121	
235/315 Fraction	1.034-1.044		1.033-1.042	
315/355 Fraction	1.098-1.125		1.094-1.122	
Xylene Insoluble, Percent				
by Weight	0.06-0.40	•	0.80-1.90	
Distillation, Percent by Weight		* V	<b>,</b>	
to 210°C	0.0-0.9		0.9-1.1	
to 235°C	3.0-10.0	• • .	1.6-6.6	
to 270°C	15.0-35.6		13.8-36.9	
to 315°C	43.4-58.0	2×1	37.6-54.2	
to 355°C	70.0-76.6		60.0-71.1	
Residue	23.7-27.5		27.4-40.0	

Source: Butala (1990)

## PRODUCTION, IMPORT, USE, AND DISPOSAL

## 4.1 PRODUCTION

Creosote refers to one of three complex mixtures of organic compounds. One mixture is derived from coal-tar (CAS Registry number 8001-58-9), one is produced from wood (e.g., beechwood) (CAS Registry number 8021-39-4), and the third is derived from the resin of the creosote bush (Larrea). Coal-tar creosote is produced from a distillation of coal-tar produced by hightemperature carbonization. Because coal-tar is a by-product of steel manufacturing, domestic production of coal-tar products is dependent on demand for steel (USITC 1987). Creosote production falls into two categories: distillate (100% creosote), and creosote in coal-tar solution. Annual production and sales figures for the years 1950 through 1987 are listed in Table 4-1. Distillate production in 1986 was 46.8 million gallons; creosote in solution was 31.6 million gallons (USITC 1987). Distillate production in 1987 was 47.3 million gallons. Production of creosote in solution in 1987 was not disclosed, but solution sales in 1987 were 34.3 million gallons (USITC 1988). The following organizations manufacture creosote domestically: Allied Corp. Morristown, New Jersey; Coopers Creek Chemical Corp, West Conshohocken, Pennsylvania; Koppers Co., Pittsburgh, Pennsylvania; Reilly Industries, Inc., Indianapolis, Indiana; Aristech Chemical Corp., Gary, Indiana; U.S. Steel Corp., Pittsburgh, Pennsylvania; Witco Chemical Corp., Woodcliff Lake, New Jersey; Crowley Tar Products Co., New York, New York; Los Angeles Chemical Co., South Gate, California; and Standard Tar Products Co., Milwaukee, Wisconsin (USITC 1988; HSDB 1988).

## 4.2 IMPORT

In 1984, 7.3 million gallons of creosote were imported into the U.S. (Bureau of Census 1984a), and 7.5 million gallons were exported from the U.S. (Bureau of Census 1984b). The is the most current data available regarding the import and export of creosote.

#### 4.3 USE

Coal-tar creosote was used as a wood preservative in the United States for over 100 years. Wood preservation accounts for over 97% of current coal-tar creosote production (Santodonato 1985). Coal-tar creosote is applied to wood by commercial pressure treatment or by home and farm dipping or brushing, although this latter use is not significant since creosote is now a restricted use pesticide (EPA 1986b). Coal-tar creosote is a wood preservative and

4. PRODUCTION, IMPORT, USE, AND DISPOSAL

Table 4-1. U.S. Production of Creosote Oila

Year		Creosote in Coal-Tar Solution <sup>b</sup> (100% creosote basis) (millions of gallons)	Distillate <sup>b</sup> (100% creosote basis) <sup>c</sup> (millions of gallons)		
1987.		47.3	(34.3)		
1986	·	46.8 (31.5)	31.6 (36.6)		
1985		64.3 (40.0)	66.6 (42.3)		
1984		42.2 (30.4)	40.8 (30.9)		
1983		39.5 (24.7)	40.9 (27.7)		
1982	•	36.3 (21.4)	44.3 (32.2)		
1981		81.0 (61.5)	61.1 (44.5)		
1980		60.6 (37.1)			
1979		80.5 (40.5)	27.2 (10.9)		
1978		51.3 (36.3)	35.0 (26.3)		
1977		47.0 (35.4)	36.0 (25.2)		
1976		77.1 (51.9)	36.8 (26.4)		
1975	.*	79.2 (50.7)	35.7 (35.2)		
1970	_	103.4	25.6		
1965		111.1	12.5		
1960		82.0	10.8		
1955	•	106.1	16.4		
1950		106.7	35.6		

<sup>&</sup>lt;sup>a</sup>Source: Santodonato 1985; USITC 1987, 1988.

bFigures in parentheses are sales figures.

Figures for 1982 and 1981 only are 100% solution basis.

# 4. PRODUCTION, IMPORT, USE, AND DISPOSAL

water-proofing agent for railroad ties, telephone poles, marine pilings, and fence posts. In addition, coal-tar creosote prevents animals and vegetable growth on concrete marine pilings, and is a component of roofing pitch, fuel oil, and lamp black and a lubricant for die molds (HSDB 1988). Other uses include animal and bird repellent, insecticide, animal dip, and fungicide (IARC 1985).

Beechwood creosote and its compounds calcium creosotate, creosote carbonate, and creosote valerate were used in the past as antiseptics and expectorants (Windholz 1983). Treatments for leprosy (Samson and Limkako 1923), pneumonia (McKinlay 1933), and tuberculosis (Fellows 1939a) also involved ingestion of beechwood creosote. Beechwood creosote is rarely used for medicinal purposes today.

#### 4.4 DISPOSAL

Creosote sludge generated from coal-tar creosote production can be fixed, solidified, and covered with clay in the settling lagoon used in treatment. This "disposal in place" requires groundwater monitoring for a 30-year period (Ball et al. 1985). Due to RCRA Land Disposal Restrictions, creosote can no longer be disposed in hazardous waste landfills unless treated to EPA-specified treatment standards (EPA 1990). Industrially used creosote-treated wood can be burned in an industrial incinerator or boiler (EPA 1986b). Treated wood used in the home or farm should be buried or disposed with household garbage, it should not be incinerated (American Wood Preservers Institute 1988).

#### 5.1 OVERVIEW

Creosote is a complex commercial mixture of some 300 organic constituents. The composition of the mixture varies from batch to batch depending on the coking process used. Creosote consists primarily of PAHs, and therefore the fate of much of the components of the mixture is similar to that of PAHs. The primary releases of creosote to the environment are in wastewater effluents from wood-treatment plants. Compared to releases to surface water and soil, creosote releases to the atmosphere are considered to be insignificant. Creosote has been found in at least 31 of the 1177 hazardous waste sites on the NPL (VIEW Database 1989). The frequency of these sites within the United States can be seen in Figure 5-1. Biotransformation by microbes is the primary process by which creosote constituents are degraded in soils, surface waters, and groundwater. The mixture is relatively stable and persistent in the environment; half-life data are not available.

Creosote has been widely used as a wood-treatment pesticide since the turn of the century. As a result of this widespread and long-term use, workers in the wood-preserving industry have been exposed to creosote for many years. Human exposure to creosote can occur by inhalation or direct dermal contact. Individuals working in wood-preserving facilities are one of the largest exposed groups. Exposure may also occur during handling and installation of treated wood products in structures such as bridges, piers, retaining walls, crossties, and fencing; as a result of burning treated scrap wood; and through contact with contaminated media at hazardous waste sites. The general public is unlikely to experience any significant exposure to liquid creosote through the direct use of wood preservative products because EPA canceled all nonwood uses of the material and restricted use of creosote products to certified applicators in January 1986 (EPA 1986b).

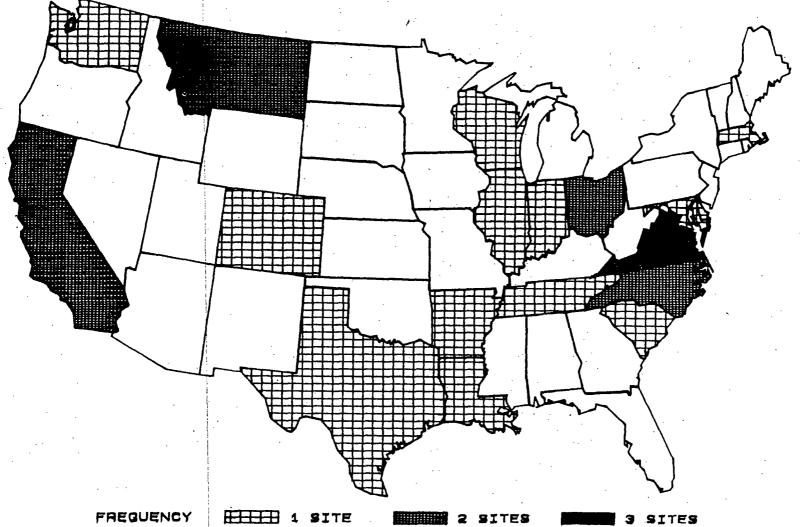
# 5.2 RELEASES TO THE ENVIRONMENT

There are no known natural sources of the creosote mixture (IARC 1973). However, several of the PAH constituents of the mixture are known to have natural sources; the reader is referred to the ATSDR Toxicological Profile for PAHs (ATSDR 1990) for additional information on natural sources, releases, and levels of PAHs not associated with creosote production, use, and disposal.

The major source of creosote released to the environment is wastewater effluents from wood treatment facilities (USDA 1980). Companies that preserve wood with creosote may treat their aqueous wastes in on-site biological treatment plants or release the wastewater into a municipal water treatment system (EPA 1978; von Rumker et al. 1975). Some creosote components may also be released to the atmosphere by fugitive emissions from these facilities.

POTENTIAL

FOR HUMAN EXPOSURE



FREQUENCY FIGURE 5-1. FREQUENCY OF SITES WITH CHEOSOTE CONTAMINATION

However, atmospheric releases are considered to be minimal in relation to releases in wastewaters.

Creosote components may also be slowly released from the surface of treated wood products by oil exudation, leaching by rainwater, or volatilization. Losses of creosote from impregnated wood are dependent on the kind of coal used to produce the coal-tar, the kind of coke oven used to make the coal-tar and the conditions under which the wood is used (Leach and Weinert 1976).

#### 5.2.1 Air

Atmospheric releases of creosote from wood-preserving plants are not well defined. Creosote constituents such as naphthalene, acenaphthylene, acenaphthene, phenanthrene, and fluorene have been detected in emissions at a pressure treatment facility that treated logs for use as utility poles and marine pilings (Engineering Science 1986). Releases may occur at several points in the treatment process, such as when cylinder doors are opened after a treatment cycle, or when creosote is transferred from the heater to the cylinder at the beginning of the impregnation process. Atmospheric releases vary from plant to plant, depending on the process design, and are considered to be significantly smaller than releases to surface water in aqueous effluents (Henningsson 1983).

On a hot, sunny day evaporation of creosote from the surface of treated wood may release creosote constituents to the atmosphere. Only the volatile creosote components such as acenaphthene and naphthalene will volatilize; the heavier fractions will remain on the wood (USDA 1980).

In a terrestrial microcosm study, release of  $^{14}$ C-labeled creosote components to the atmosphere from treated wood accounted for 1.0% of total acenaphthene and 1.4% phenanthrene whereas 93.5% and 95% of these components, respectively were retained in the wood (Gile et al. 1982).

Other potential sources of atmospheric releases include incineration of scrap wood treated with the mixture and re-entrainment of dust and soils contaminated with components of the mixture in the vicinity of hazardous waste sites.

### 5.2.2 Water

The major source of creosote released into surface waters and groundwater is wastewater effluents from wood-preserving facilities (USDA 1980). In previous years, wastewater generated from wood treatment facilities was often discharged to unlined evaporation/settling lagoons where a sludge was formed. Water-soluble creosote components then percolated through the soil to reach the groundwater table. Wastewaters may include process water generated from steam conditioning of the wood, preservative formulation recovery and

regeneration water, water used to wash excess preservative from the surface of the wood, condensate from drying kilns used to dry preserved or surface, protected wood, water that accumulates in door and retort sumps, and rain falling on or in the immediate vicinity of the treating cylinder and work tank area. Groundwater contamination from creosote wastewaters and sludge stored in unlined surface water impoundments at a wood treatment facility has been reported in Pensacola, Florida (Goerlitz et al. 1985). Similar contamination problems have been reported in Conroe, Texas (Borden 1986), and St. Louis Park, Minnesota (Hickock et al. 1982).

Water-soluble creosote constituents (e.g., phenols) may be released to surface water or groundwater by leaching from the surface of creosote contaminated soils at hazardous waste sites or from treated wood products coming into contact with water, such as marine pilings. For example, some studies have shown that creosote is lost to a greater extent from marine timber than from timber placed in freshwater as a result of cell contraction caused by the high concentration of salts in seawater (Henningsson 1983).

#### 5.2.3 Soil

Creosote may be released to soils at treatment facilities as a result of bleeding of the product from treated timber in stock yard and storage areas and through disposal of the mixture at hazardous waste sites. Rain water may also wash the soluble components directly from the surface of treated timber and into the soil (Henningsson 1983).

## 5.3 ENVIRONMENTAL FATE

# 5.3.1 Transport and Partitioning

Creosote constituents released to surface waters will differentially partition to the water column or to sediments depending on their water solubility and sorptive properties. For example, PAHs, the major constituents of creosote, generally tend to sorb strongly to soil and sediment particulates and have low aqueous solubilities and mobility (Hickock et al. 1982).

Nitrogenous bases present in creosote wastewater (e.g., aniline, toluidines, and xylidines) are relatively soluble, mobile, and persistent in groundwater (Pereira et al. 1983). However, behavior at a given site is also dependent on site-specific characteristics. For example, PAH, phenol, and heterocyclic components of creosote wood treatment process wastes were found to migrate en masse in groundwater through a contaminated sand and gravel aquifer in Pensacola, Florida; sorption of these different classes of organic constituents in the low organic carbon(<0.1%) aquifer materials was not important (Pereira and Rostad 1986).

In an investigation of the extent of creosote contamination at four wood preservative plants with process water surface impoundments, unspecified creosote components were found to have moved 20-60 feet vertically from the

impoundments to the water table and up to 500 feet horizontally from the sources (Ball 1987).

In an investigation of the release of creosote from treated wood into freshwater and sea water, naphthalene, phenanthrene, acenaphthene, dibenzofuran, fluorene, and 2-methylnaphthalene were found to be the major components that migrated into water (Ingram et al. 1982). The rate of migration was found to increase significantly with increasing temperature within the range of 20° to 40°C and occurred more slowly from aged than from freshly treated pilings.

In a terrestrial microcosm study, 2.7% of radiolabeled phenanthrene and 4.3% of radiolabeled acenaphthene were found in soil samples taken in a 10-cm zone around creosote-treated posts, whereas concentrations of the compounds in the posts were 95% and 93.5% of the amounts applied, respectively, after 2.5 months (Gile et al. 1982).

Limited uptake of some creosote constituents has been detected in plants exposed to creosote-treated wood in nearby soil. Only 0.04% of applied acenaphthene and 0.1% of phenanthrene partitioned to plants in one study (Gile et al. 1982).

Animals such as voles, crickets, snails, pill bugs, and worms have exhibited the capacity to assimilate radiolabeled creosote components in terrestrial microcosm studies. Creosote components were found to accumulate to the greatest extent in the vole, with bioconcentration factors of 12-31. The <sup>14</sup>C mass balance content of the animals was 1.2% of applied acenaphthene and 0.8% of applied phenanthrene versus 4.3% and 2.7%, respectively, in soils (Gile et al. 1982). In addition, mussels taken from creosote-treated pilings have been found to contain significantly more B[a]P, a creosote constituent, than those growing elsewhere (Dunn and Stich 1976). Accumulation of creosote-derived PAHs has been reported in benthic organisms in Pensacola Bay (Elder and Dresler 1988; Rostad and Pereira 1987). Fluoranthene, pyrene, B[a]P, anthracene, chrysene, and phenanthrene were detected in higher concentrations in tissues of snails (Thais haemastoma) and oysters (Crassostrea virginica) taken from offshore sites near an onshore wood treatment plant than from control sites.

### 5.3.2 Transformation and Degradation

#### 5.3.2.1 Air

No information was found in the available literature concerning the transformation of creosote components in the atmosphere.

#### 5.3.2.2 Water

Creosote components are degraded in aquatic environments by microfaunal metabolism (Borthwick and Patrick 1982; Ingram et al. 1982). Microorganisms may act on the creosote-treated wood itself or on creosote components that have leached from the treated wood. Quinoline, the major tar base in creosote, has been reported to be degraded in surface water and groundwater by bacteria of the genus <u>Pseudomonas</u> (Bennett et al. 1985). Biotransformation of the phenolic components of creosote apparently also occurs under anaerobic conditions in contaminated groundwater (Ehrlich et al. 1983; Goerlitz et al. 1985). Adaptation of soil microorganisms to PAH contaminants in groundwater originating from creosote treatment plant wastes has also been reported (Wilson et al. 1986).

Creosote components have been detected in surface water samples taken near a wood-treatment facility that ceased operation 30 years earlier (Black 1982). The creosote appeared to have been transported through the soil and was entering the river via seepages and springs. Weathering processes produced only minor constitutive changes in the creosote with relative losses of the lower molecular weight components. These changes probably reflected the greater volatility and solubilities of the two and three carbon ring PAHs.

## 5.3.2.3 Soil

Creosote components are slowly released from treated wood products by oil exudation, rainwater leaching, and by volatilization of the lighter fractions (Henningsson 1983). USDA (1980) reported that the major components of creosote were not detected in soil samples taken to a depth of 6 inches within 2-24 inches from treated poles, presumably as a result of biotransformation of mobilized components by soil microorganisms. Creosote components released to soils in wastewater effluents have been found to be biotransformed by soil microbes under aerobic conditions (Middleton 1984). Bacteria of the genus Pseudomonas isolated from a creosote-contaminated waste site have been reported to degrade creosote-derived quinoline (Bennett et al. 1985). Acclimation to creosote phenolic constituents by soil microorganisms has also been demonstrated (Smith et al. 1985).

## 5.4 LEVELS MONITORED OR ESTIMATED IN THE ENVIRONMENT

#### 5.4.1 Air

No information was found in the available literature regarding ambient atmospheric concentrations of creosote-derived components (i.e., PAHs) in the United States. Workplace air concentration data are discussed in Section 5.5. Data on ambient atmospheric concentrations of PAHs derived from other sources can be found in the ATSDR Toxicological Profile for PAHs (ATSDR 1990).

#### 5.4.2 Water

Results from 2 years of groundwater sampling at an abandoned wood treatment facility in Conroe, Texas where creosote had been used for about 20 years showed that wells were contaminated with levels of up to 3490  $\mu$ g/L naphthalene, 1263  $\mu$ g/L methylnaphthalene, 425  $\mu$ g/L dibenzofuran, and 302  $\mu$ g/L fluorene. The contaminants had apparently migrated through the clay and sand soils on the site from three waste pits (Bedient et al. 1984).

### 5.4.3 Soil

Several PAH constituents of creosote were detected in soil samples taken at an abandoned wood treatment facility in Conroe, Texas at depths of up to 25 feet. Maximum concentrations of the compounds were detected in samples collected at the 0.7-1.8 ft depth. Maximum concentration levels were 3.7 mg/kg for naphthalene, 3.4 mg/kg for methylnaphthalene, 3.8 mg/kg for dibenzofuran, 4.2 mg/kg for fluorene, and 2.2 mg/kg for anthracene (Bedient et al. 1984).

### 5.4.4 Other Media

Creosote-derived phenanthrene, 1,2-benzanthracene, and B[a]P have been detected in river sediments at concentrations of up to 231,000  $\mu$ g/kg, 62,000  $\mu$ g/kg, and 16,000  $\mu$ g/kg, respectively, downstream from the site of a former wood treatment facility (Black 1982). Creosote-derived PAHs were also detected in the sediments of Pensacola Bay and a drainage stream in the vicinity of a former wood treatment facility near Pensacola, Florida. PAH concentrations ranged from 200  $\mu$ g/g for naphthalene to 140,000  $\mu$ g/kg for anthracene in stream sediments; concentrations in Pensacola Bay ranged from 75  $\mu$ g/kg for benzanthracene to 190  $\mu$ g/kg for fluoranthene (Elder and Dresler 1988).

#### 5.5 GENERAL POPULATION AND OCCUPATIONAL EXPOSURE

Individuals working in the wood-preserving industry comprise the largest portion of the population potentially exposed to creosote. Workers employed at creosote pressure treatment facilities may be exposed by direct dermal contact or by inhalation of volatilized components. Approximately 4,000 workers in the United States were employed at 188 pressure treatment plants that used creosote as a preservative in the late 1970s (USDA 1980). Potential exposure to creosote in these plants is minimized by the use of closed systems for receiving, transferring, mixing, storing, and applying the mixture to wood products. Similarly, dermal exposure from the handling of freshly treated wood is minimized by the use of highly mechanized processes. Exposure via inhalation, however, is more likely to occur. For example, worker exposure may be significant during opening of treatment cylinder doors and cylinder cleaning operations (EPA 1981b). Inhalation and dermal exposure are also more likely in plants using nonpressure treatment methods such as thermal and dip

treatments in open tanks. An estimated 100 workers were involved in commercial thermal and dip treatment operations in the late 1970s. Some of these workers experienced consistently high inhalation exposures (USDA 1980). Other historical nonpressure treatment exposures included an estimated 50,000 individuals (e.g., homeowners, farmers, landscapers) who applied creosote in noncommercial brush, dip, spray, and soak treatments (EPA 1981a). Dermal contact and inhalation may have resulted in exposure to high concentrations of creosote components for these individuals, but the exposures were usually of intermittent frequency (USDA 1980). However, designation of creosote products as restricted use pesticides by EPA in 1986 has probably decreased the number of individuals potentially exposed in these nonpressure wood treatment applications (EPA 1986b).

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There are currently 25,000 workers employed in about 75-100 domestic wood treatment plants using creosote. As a result of the use of engineering controls and personal protective equipment (e.g., respiratory protection and impervious gloves) required in the 1986 settlement of the EPA Special Review process<sup>2</sup>, airborne exposures to creosote components in the workplace are generally below the OSHA permissible exposure limit (PEL) of 0.2 mg benzene soluble particulates/m³ air (Rivers 1990).

However, prior to the standard use of these controls by industry, workers were potentially exposed to higher airborne concentrations of creosote constituents. For example, the concentration of creosote (i.e., coal-tar pitch volatiles) components in personal air samples taken at a railroad tie treatment facility in Somerville, Texas, were found to range from 0.003 to 1.211 mg/m³ (Todd and Timbie 1980). Another industrial hygiene survey of worker exposure to creosote at a wood-treatment facility in Tacoma, Washington, showed coal-tar pitch volatiles in personal air samples ranged from less than 0.4  $\mu$ g/m³ to 111.9  $\mu$ g/m³ (Todd and Timbie 1981). The higher concentrations were found at the end of the treatment process when the cylinder was opened. NIOSH investigated creosote exposure among dock builders in Brooklyn, New York, in 1980. Employees were reported to have substantial direct skin contact with creosote. Breathing zone concentrations of the cyclohexane extractable fraction of the coal-tar pitch volatiles ranged from zero to 0.059 mg/m³ of air (Baker and Fannick 1981). Monitoring data

<sup>&</sup>lt;sup>2</sup>A Special Review of a currently registered pesticide may be initiated by EPA when validated data indicate that certain types of toxicity (e.g., carcinogenicity, developmental toxicity, acute effects) exist for humans or for non-target plant or animal species. A formal process exists for notifying registrants and other interested parties, requesting further data regarding the pesticide in question, analyzing and reporting risks and benefits, and requesting public review. The final regulatory decision may be implemented over a period of time or it may be imposed immediately as an emergency action based solely on the EPA's finding of immediate danger to human health or the environment.

collected from 11 creosote treatment facilities in the United States showed that worker breathing zone concentrations of benzene-soluble particulates averaged 0.07 mg/m³ (range 0.01-0.061 mg/m³) (EPA 1981a). Comprehensive studies of worker exposure to creosote in wood treatment plants have been conducted by Koppers Company alone and in conjunction with NIOSH (SRI International 1977; Markel et al. 1977). Data from these studies indicated that, on the average, employee exposure to particulate polycyclic organic materials (PPOM) was within the permissible level of 0.1 mg/m³ recommended by NIOSH for coal-tar volatiles. The only components that could be reliably measured in the vapor-phase fractions collected were naphthalene, methylnaphthalene, and acenaphthene. The concentrations of these chemicals ranged from 0.54 to 2.0 mg/m³. Fluorene and phenanthrene-anthracene were detected in trace quantities but were not quantifiable. Benzene-soluble particulates (PPOM) ranged from 0.02 to 0.10 mg/m³.

A gravimetric analytical method has been used in most of these workplace monitoring studies. This method involves the collection of airborne particulates on glass fiber filters and subsequent extraction by solvents, such as benzene or cyclohexane. The extracted fraction of the particulate matter is determined by weighing. As a result of two significant shortcomings of this method, the inability to identify constituents of the airborne particulates and to sample vapor phase components, EPA (1981a) concluded that definitive information was not available on the identity of airborne components of creosote in workplace atmospheres. EPA (1981a) also stated that quantitative estimates of treatment plant worker dermal exposures and quantitative inhalation or dermal exposure data for workers applying creosote in nonpressure treatment scenarios and for downstream workers who install, handle, or contact treated wood products were unavailable. Vapor phase components were found to be an important source of worker exposure in an industrial hygiene survey of wood treatment facilities conducted in Finland. Most of the airborne contaminants in worker breathing zones were in the vapor phase; the proportion of particulate PAHs to total concentration of vapors was <0.5%-3.7% (Heikkila et al. 1987).

Exposure to creosote may also occur during installation of treated poles, during inspection and maintenance operations, and through casual contact (USDA 1980). One of the major end point uses of creosote is treatment of railroad crossties. Since crossties are installed mechanically by railroad companies, workers generally have minimal dermal exposure in this process. Exposure via inhalation, however, is considered to be moderate and consistent during this type of installation procedure. In other situations, crossties may be installed manually, in which case there is consistent moderate to high exposure via skin contact as well as by inhalation. The amount of exposure via skin contact ranges from moderate to high depending on whether workers wear protective clothing. Skin contact is considered minimal for railroad personnel who inspect ties in service as well as for the general public who may have casual contact with creosote-treated crossties. In instances where crossties are used for landscaping purposes, contractors involved in the sale

and installation of freshly treated ties experience consistent moderate exposure via inhalation and minimal to occasionally high exposure for skin contact.

Installation of treated lumber and timbers in structures such as bridges, piers, retaining walls, fences, and barns involves a significant amount of manual contact. Likewise, the installation of switch ties, cross planks, crossarms, block flooring, and fence posts is usually done manually. In these situations, human exposure via inhalation is considered moderate while exposure via skin contact may vary from minimal to high depending on the type of protective equipment used (USDA 1980).

The opportunity for direct dermal contact is particularly high for workers installing treated poles. Activities such as attaching fittings often preclude the use of protective gloves, and as a result of creosote bleeding from the treated poles, the potential for dermal contact of workers performing maintenance operations persists for years after installation (Henningsson 1983).

Exposure of individuals installing treated fence posts and lumber and timbers via inhalation of creosote volatiles (e.g., acenaphthene and naphthalene) can also occur when freshly treated materials are handled under calm, het, sunny conditions (USDA 1980).

Potential sources of non-occupational human exposure to creosote include contact with creosote-treated wood products (e.g., railroad ties used for landscaping), incineration of creosote-treated scrap lumber, and contact with contaminated environmental media at hazardous waste sites (e.g., ingestion of contaminated groundwater). Direct exposure of homeowners to wood treatment products containing creosote should be limited, since EPA has restricted the sale and use of such products to certified applicators. Industrial sources have noted that there have been no reports or instances of health effects allegations, except for rare reports of skin irritation, resulting from public contact with creosote-treated wood in the last 20 years.

The only information on biological indicators of exposure to creosote found in the available literature involved a study of 1-hydroxypyrene in the urine of a creosote wood treatment plant worker (Jongeneelen et al. 1985).

## 5.6 POPULATIONS WITH POTENTIALLY HIGH EXPOSURES

Individuals living in the vicinity of hazardous waste sites and abandoned wood treatment plants contaminated with creosote may experience higher levels of exposure than the rest of the general population.

Individuals who apply creosote directly to wood, including farmers, carpenters, and homeowners who come in contact with creosote-treated wood products are believed to be exposed to the highest levels of creosote

components via inhalation and dermal contact. It has been estimated that historically about 4,000 workers may have been routinely exposed and up to 50,000 people may have been intermittently exposed to creosote through its application as a preservative to wood products (USDA 1980). The size of this population has probably decreased over the last few years since EPA restricted the use of creosote to certified applicators.

# 5.7 ADEQUACY OF THE DATABASE

Section 104(i) of CERCLA directs the Administrator of ATSDR (in consultation with the Administrator of EPA and agencies and programs of the Public Health Service) to assess whether adequate information on the health effects of creosote is available. Where adequate information is not available, ATSDR, in conjunction with the National Toxicology Program (NTP), is required to assure the initiation of a program of research designed to determine the health effects (and techniques for developing methods to determine such health effects) of creosote.

The following categories of possible data needs have been identified by a joint team of scientists from ATSDR, NTP, and EPA. They are defined as substance-specific informational needs that, if met would reduce or eliminate the uncertainties of human health assessment. In the future, the identified data needs will be evaluated and prioritized, and a substance-specific research agenda will be proposed.

## 5.7.1 Identification of Data Needs

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Physical and Chemical Properties. Limited physical property data, such as boiling point and density (see Table 3-2), are available for the coal-tar creosote mixture. Additional physical and chemical property data, such as water solubility, vapor pressure,  $K_{\rm oc}$ , and Henry's law constant values would be useful in order to predict the partitioning and transformation of coal-tar creosote components in air, water, and soil. These values are currently not available because their determination is complicated by the fact that creosote is a mixture of variable composition.

Production, Use, Release, and Disposal. Manufacturing methods are well described in the literature. Production figures are limited because of the confidential nature of this business information. Uses of creosote, both coal-tar and beechwood, are well described. Since the use of coal-tar creosote as a wood preservative has been restricted, the potential of the population to be exposed is greatly diminished. The major releases of creosote resulting from treatment processes at wood-preserving plants are known but the levels are not well quantified. Current production, release, and disposal information would assist in identifying the levels of creosote present in the environment and, thus, populations potentially exposed as a result of these processes. Creosote sludge from production processes can be

treated and disposed on-site with proper groundwater monitoring. Creosote can no longer be disposed in hazardous waste landfills unless treated to EPA-specified standards. Creosote-treated wood used in industrial applicators can be burned in an industrial incinerator or boiler; however, treated wood used in domestic or farm applications should be buried rather than incinerated.

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Environmental Fate. The limited information available regarding transport and partitioning of creosote components among environmental compartments indicates mobility of PAH, phenol, and heterocyclic constituents of the mixture in water, sorption of PAH components in soils, and bioconcentration of creosote-derived PAHs by terrestrial and aquatic organisms. Biotransformation appears to be the most important degradation process in soils and aquatic environments. Additional data on the transport of volatile creosote components in the atmosphere and the partitioning of creosote released to surface waters and soils would be useful. Quantitative data on the rates of biotransformation in soils, surface water, and groundwater under aerobic and anaerobic conditions would also be useful. importance of other transformation processes, such as photolysis, photooxidation, and hydrolysis, in relation to biotransformation and rates of transport between media, should also be defined. These data would be useful to help define potential pathways of human exposure and to estimate ambient concentrations of creosote components in environmental media.

Bioavailability from Environmental Media. Very limited information was found in the available literature regarding the uptake of creosote components by living organisms from contaminated water and soil at hazardous waste sites. Studies have been done with persistent constituents (e.g., PAHs) which show that plant uptake from soils is limited, whereas bioconcentration in aquatic organisms from contaminated surface waters has been demonstrated. Data from human and animal studies indicate that creosote components are absorbed following ingestion, inhalation, and dermal exposure to the mixture. PAH and phenol constituents have also been demonstrated to be absorbed following exposure by these routes. However, quantitative information regarding the uptake and bioavailability of creosote from contaminated environmental media is not available. In addition, data are needed on the bioavailability of creosote-contaminated soils.

Food Chain Bioaccumulation. Very limited information was found in the available literature regarding the biomagnification of creosote-derived compounds among food chain trophic levels. However, PAHs, the major persistent constituents of the commercial mixture, have not been reported to undergo significant biomagnification in aquatic food chains because of the ability of many aquatic organisms to rapidly metabolize and eliminate these compounds (Eisler 1987). Since components of the mixture are not concentrated to high levels in human food items, food chain bioaccumulation does not appear to be an important source of human exposure and no additional information is needed at this time.

Exposure Levels in Environmental Media. Limited information is available regarding ambient concentrations of creosote-derived PAHs in soils and no data are available regarding atmospheric concentrations of creosote components. Very limited information is available on concentrations of component compounds in surface waters and sediments receiving wood treatment plant effluents. Data are particularly lacking for contaminated media in the vicinity of hazardous waste sites. These data would be useful to estimate the exposure of populations coming into contact with components of the mixture through inhalation of contaminated air, consumption of contaminated surface water or groundwater, or direct dermal contact with environmental media.

Exposure Levels in Humans. A population exists that is potentially exposed to creosote through contact with contaminated media at hazardous waste sites and with treated wood products. A second potentially exposed workforce population exists at wood treatment facilities. Currently no information exists that demonstrates tissue levels of any components of the mixture in these populations. Estimates of human exposure to creosote constituents, or body burdens of creosote components, are complicated by the lack of information on exposure to creosote constituents and levels of creosote-derived components in the environment. Collecting information on tissue levels of creosote components in humans would be necessary to examine the relationship between levels of creosote-derived compounds in the environment, human tissue levels, and subsequent development of health effects.

Exposure Registries. No exposure registries for creosote were located. This compound is not currently one of the compounds for which a subregistry has been established in the ATSDR National Exposure Registry. The compound will be considered in the future when chemical selection is made for subregistries to be established. The information that is amassed in the ATSDR National Exposure Registry facilitates the epidemiological research needed to assess adverse health outcomes that may be related to the exposure to this compound.

#### 5.7.2 On-going Studies

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Creosote is currently subject to a FIFRA registration standard and data call-in by EPA. In addition, the Creosote Council II is planning to conduct a research program that includes testing for worker exposure and protection.

Remedial investigation/feasibility studies being conducted at the 31 NPL sites where creosote has been found and at the numerous creosote-contaminated RCRA corrective action sites should provide data on concentrations of the mixture in contaminated media in the vicinity of hazardous waste sites. For example, creosote constituents have been found in surface water impoundments and soil samples taken near wood treatment facility sites in Colorado, Louisiana, Texas, and Montana.

#### 6. ANALYTICAL METHODS

The purpose of this chapter is to describe the analytical methods that are available for detecting and/or measuring and monitoring crossote mixtures in environmental media and in biological samples. The intent is not to provide an exhaustive list of analytical methods that could be used to detect and quantify the components of crossote. Rather, the intention is to identify well-established methods that are used as the standard methods of analyses. Many of the analytical methods used to detect crossote mixtures in environmental samples are the methods approved by federal agencies such as EPA and the National Institute for Occupational Safety and Health (NIOSH). Other methods presented in this chapter are those that are approved by a trade association such as the Association of Official Analytical Chemists (AOAC) and the American Public Health Association (APHA). Additionally, analytical methods are included that refine previously used methods to obtain lower detection limits, and/or to improve accuracy and precision.

The analytical methods used to quantify creosote mixtures in biological and environmental samples are summarized below. Table 6-1 lists the applicable analytical methods for determining components of creosote in biological fluids and tissues and Table 6-2 lists the applicable analytical methods for determining the components of creosote in environmental samples.

## 6.1 BIOLOGICAL MATERIALS

Coal-tar creosote contains a highly complex mixture of organic components. These include PAHs (polycyclic aromatic hydrocarbons), phenols, sulfur, oxygen and nitrogen heterocycles. Various PAHs have been found to constitute about 85% of coal-tar creosote (Bedient et al. 1984). The levels of creosote in biological materials can therefore be estimated by measuring the PAH content in biological samples. The methods available to measure PAHs in biological materials include gas chromatography equipped with a flame ionization detection (GC/FID), gas chromatography coupled to a mass spectrometry (GC/MS) and high performance liquid chromatography (HPLC). Immunoassay techniques, i.e., enzyme linked immunosorbent assays (ELISA) and ultrasensitive enzyme radio immunoassay (USERIA), <sup>32</sup>P-postlabeling and synchronous luminesence spectroscopy (SLS) are methods currently being developed to detect and quantify ultratrace levels of PAH adducts bound covalently to macromolecules (e.g., DNA).

GC/MS and HPLC has been employed to detect creosote-derived PAH complexes at ppt (pg/g) levels in human tissues (Liao et al. 1988; Obana et al. 1981). The detection and quantification of trace levels of PAHs in biological tissues involves extensive and rigorous clean-up procedures including florisil, silica and alumina column chromatography (Liao et al. 1988; Obana et al. 1981).

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TABLE 6-1. Analytical Methods for Determining Croosote-Derived PAH Components in Biological Materials

Sample Matrix	Sample Preparation	Analytical Method	Sample Detection Limit	Accuracy	Reference
Adipose tissues	Add 8% benzene in hexane to adipose tis- sue sample and homogenize; add Na <sub>2</sub> O <sub>4</sub> ,	GC/MS	5-50 ng/g	521-951 recovery	Liso et al. 1988
	stir and load onto florisil column; elute column with 8% benzene in hexane, concentrate, and analyze		• · · · · · · · · · · · · · · · · · · ·	. *	:
iver	Add S-9 mix (KH <sub>2</sub> PO <sub>4</sub> buffer, MgCl <sub>2</sub> , KCl, glucose-6-phosphate, NADF*) at pH 7.4 to homogenized sample in DMSO and incubate	RPLC	No data	No data	Amin et al. 1982
	with shaking at 37°C; add ice-cold ace- tone to stop reaction and extract with ethyl acetate; dry (Na <sub>2</sub> SO <sub>4</sub> ), concentrate,	!			
	and analyze metabolites  Mince tissue sample, saponify with	HPLC	0.006-0.46	No data	Obana et al. 1981
	KOH/EtoH, and extract with hexane; separate hexane layer and extract with DMSO; pour DMSO layer into 15% aqueous		ng/g		:
	MaCl solution and reextract with hexane; wash hexane layer with water, dry (Na_SO_s), and concentrate; load extract	•			• •
	onto silica or alumina gel column; elute column with hexane followed by diethyl ether:hexane (15:85) and concentrate				
	eluent				
lood	Collect blood sample in heparinized plastic tube and centrifuge; separate white blood cells and isolate DNA by standard RNase and phenol treatment;	ELISA	1x10 <sup>-15</sup> mol BPDE per µg DNA	Ro data	Perera et al. 1986
	assay BPDE-NDA adduct by immunoassay		• .	•	
	Isolate PAR-DNA adduct from white blood cells of blood; digest adduct with radiolabeled (32P)ATP; resolve the	32P-Postlab- eling	0.3x10 <sup>-15</sup> mol adduct per µg DNA	No data	Philips et al. 1988
	radiolabeled adducts on TLC plate and detect by counting radioactivity				
	Collect blood sample in heparinized plastic tube and centrifuge; collect lymphocyte cells and isolate BPDE-DNA adduct by standard treatment; analyze	ELISA or USERIA and SLS	0.006- 0.23x10 <sup>-15</sup> mol BPDE per µg DNA	No data	Harris et al. 1985

TABLE 6-1 (Continued)

Sample Matrix	Sample Preparation	Analytical Method	Sample Detection Limit	Accuracy	Reference
Urine	Inject animal with radiolabeled benzo(a)pyrene; collect urine sample, add MeOH and load onto C-18 Sep-Pak Column; elute with aqueous MeOH and analyze eluent by HPLC	HPLC	5x10 <sup>-12</sup> mol 7-BPDE- Gua per μg of labeled benzo(a) pyrene	No data	Autrup and Seremet 1986

GC/MS = gas chromatography/mess spectrometry; HPLC = high performance liquid chromatography; ELISA = enzyme linked immunosorbent assay; USERIA = ultra-sensitive enzyme radioimmunoassay; SLS = synchronous luminescence spectroscopy; NADP\* = oxidized nicotinamide adenosine dinucleotide; DMSO = dimethyl sulfoxide; BPDE = benzo(a)pyrene diol epoxide; and Gua = guanine.

TABLE 6-2. Analytical Methods for Determining Creosote-Derived PAH Components in Environmental Samples

Sample Matrix	Sample Preparation	Analytical Method	Sample Detection Limit	Accuracy	Reference	
. Wooden sleepers (railroad	Extract accurately weighed sample with diethylether; filter extract through anhydrous sodium sulfate and evaporate	GC/MS	1-3 ng/ sample	No data	Rotard and Mailahn 1987	
crossties) in play ground	the solvent; perform acid/base/neutral liquid-liquid separations on creosote residue			·		
Coal-tar creosote	Dissolve sample in cyclohexane and ex- tract with 90% methanol; evaporate ex- tract to dryness; dissolve residue with cyclohexane and extract with nitrome- thane; evaporate extract to dryness and	<b>GC</b>	10 ppm	No data	Lijinsky et al. 1963	6.
	dissolve residue with small amount of benzene  Dissolve sample in methylene chloride at a concentration of ~10% (W/w); inject sample solution into g.c. for analysis	сс	No data	No data	Nestler 1974	ANALYTICAL
River sedi- ments	Digest wet sediment sample in hoiling EtOH/KOH; partitioned hydrocarbons by extraction into cyclohexane; concentrate extract and load onto florisil column; elute PAH complex from column with 50% methylene chloride/hexane; concentrate sample and analyze	HPLC	No data	No data	Black 1982	CAL METHODS
Contaminated groundwater	Filter sample through prebaked glass- fiber filters to remove suspended sedi- ments; pass sample through a bonded-phase extraction column; elute trapped organic compound from column with acetonitrile followed by methylene chloride; dry elute with Na <sub>2</sub> SO <sub>4</sub> and concentrate with a stream of dry nitrogen	GC/MS	50 μg/L	95% recovery	Rostad et al. 1984	
Impregnated wood (work- place)	Heat sample at 60.C in a chamber and load onto an XAD-2 resin column; extract resin with diethyl ether and analyze	GC/MS	0.01- 0.005 mg/m <sup>3</sup>	82%-102% re- covery	Heikkila et al. 1987	

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TABLE 6-2 (Continued)

Sample Matrix	Sample Preparation	Analytical Method	Sample Detection Limit	Accuracy	Reference
	Collect heated sample on a prewashed (cy-	HPLC	8 ng/m <sup>3</sup>	No data	Heikkila et al. 1987
	clohexane) glass fiber filter; extract sample with cyclohexane and evaporate to dryness; dissolve residue in acetoni-				
	trile: water (85:15)		,		. :
Creosote treated wood	Place a small piece of treated wood in a solid sample holder in the heated injection port of G.C.; vaporize the creosote from wood onto column and analyze	GC	No data	No data	Lorenz and Gjovik 1972
Gas and par- ticulate	Pump sample through a glass fiber filter- amberlite XAD-2 adsorbent sampling sys-	HPLC	0.005 to 2.5 mg/m <sup>3</sup>	871-1021 re- covery	Anderson et al. 1983
matter (workplace)	tem; extract with diethylether in an ultrasonic bath; concentrate ether extract and dilute with acetonitrile and analyze				

GC/MS = gas chromatography/mass spectrometry; GC = gas chromatography; HPLC = high-performance liquid chromatography; and FAH = polycyclic aromatic hydrocarbon.

There is considerable evidence that PAHs are enzymatically converted to highly reactive metabolites that bind covalently to macromolecules such as DNA, thereby causing carcinogenesis and mutagenesis in mammalian systems. Thus, benzo(a)pyrene a prototype of the carcinogenic PAHs and the most thoroughly studied PAH, is converted by specific cellular enzymes to the synand anti-isomers of  $7\beta$ ,  $8\delta$ -dihydroxy-( $9\delta$ ,  $10\delta$ )-epoxy-7, 8, 9, 10-tetrahydrobenzo(a)pyrene (BPDE) and binds covalently to DNA, resulting in formation of the putative BPDE-DNA adduct (Harris et al. 1985; Haugen et al. 1986; Autrup and Seremet 1986).

The ELISA technique has been employed for detecting antibodies in serum bound to BPDE-DNA adducts. The USERIA method involves measuring the immunological response of BPDE-DNA in the presence of rabbit anti-serum, alkaline phosphatase enzyme and radiolabeled para-nitrophenyl phosphate (PNPP). The radioactivity of the hydrolysed tritiated PNPP is measured by a scintillation counter. Both ELISA and USERIA methods has been employed to detect PAH-DNA adducts at  $10^{-15}$  mol levels in the blood and tissues of humans occupationally exposed to PAH (Amin 1982; Harris et al. 1985; Haugen et al. 1986; Newman et al. 1988; Perera et al. 1988). The 32P-postlabeling method involves a 5'-labeling of DNA-adducts that has been digested with nuclease P1 enzyme system to 3'-mononucleotides. Adducts present in the digest that were resistant to nuclease P1 were thus labeled with 32P, while unmodified nucleotides were not. The digested DNA adducts are separated by thin-layer chromatography and quantified by scintillation counting. A detection limit of  $0.3 \times 10^{-15}$  mol of PAH adduct per  $\mu g$  of DNA (less than one adduct in  $10^7$ nucleotides) has been achieved (Philips et al. 1988).

### 6.2 ENVIRONMENTAL SAMPLES

Coal-tar creosote is a amber to brown oily liquid obtained by the fractional distillation (at 200°-400°C) of crude coal-tars. The commercial product is a highly complex mixture of hundreds of components, including PAHs, phenols, and nitrogen, oxygen, and sulfur heterocycles. The mixture per se cannot be measured. The PAH component fraction is most often used as an indicator of creosote contamination of environmental media. For example, screening for total PAHs is often used at hazardous waste sites when creosote contamination is suspected. The PAH fraction is used in these analyses because it is more persistent than the phenolic fraction, which tends to be more mobile and biodegradable. The methods used to measure total PAHs also detect the nitrogen, oxygen, and sulfur heterocyclic components of the mixture.

The determination of trace amounts of creosote-derived PAHs in environmental samples has been restricted to a limited number of techniques. These include high-performance liquid chromatography (HPLC), gas chromatography equipped with flame ionization detection (GC/FID), and gas chromatography coupled to chemical ionized mass spectrometer (GC/MS).

GC/FID or GC/MS are perhaps the most widely employed analytical techniques for the determination of creosote-derived PAHs in coal-tar, contaminated groundwater, wooden sleepers (railroad crossties), and impregnated wood (Heikkila et al. 1987; Lijinsky et al. 1963; Lorenz and Gjovik 1972; Nestler 1974; Rotard and Mailahn 1987; Rostad et al. 1984). GC/FID is one of the methods recommended by EPA for detection of PAHs in wastewater and solid waste (EPA 1986c).

Heikkila et al. (1987), employed GC/MS technique to determine creosote levels of impregnated wood in the workplace. Detection limits of  $10 \times 10^{-6}$  to 50x10<sup>-6</sup> g of creosote per m<sup>3</sup> of sample and recoveries of 82% and 102% were achieved. Heikkila and co-workers measured the components of PAHs with a reverse phase HPLC equipped with fluorescence detection. For the detection of creosote vapors, naphthalene was used as an indicator since it constitutes about 18% by weight of total PAHs in creosote (Andersson et al. 1983; Heikkila et al. 1987). Rotard and Mailahn (1987) used a modified sample extraction procedure to identify various components of creosote extracts in wooden sleepers. The procedure involved the separation of compounds by functional group using acid, base, and neutral conditions. Detected compounds include phenanthrene, anthracene, and naphthalene (neutral extractions), quinoline and isoquinoline (basic extraction), cresols, and phenols (acidic extraction). Rostad and co-workers (1984) developed a method for isolation and detection of creosote in contaminated groundwater. This method involves passing the sample through a small column containing a solid-bonded phase adsorbent, which adsorbs the organic compounds. The authors indicated that this method is simple, faster, and cheaper to perform than the acid/base/neutral extraction procedure. It effectively isolated all organic compounds from contaminated groundwater regardless of polarity, functional group, or water solubility in one step, thereby minimizing hazardous exposure to sample.

HPLC equipped with fluorescence detection has been used to identify creosote-derived PAH complex in river sediments (Black 1982). Complex peaks of PAH compounds were observed using an excited wavelength at 300 nm and an emitted wavelength at 420 nm. Some PAHs, such as phenanthrene and B[a]P, were completely resolved using appropriate chromatographic conditions. Benz[a]anthracene was only partially resolved (absorbance at 254 nm) from a complex of peaks associated with the chrysene-benzene[a]anthracene doublet (Black 1982). Andersson et al. (1983) employed an amberlite XAD-2 adsorbent instead of filter system for isolating organic compounds from gas and particulate matter in creosote impregnating plant. Good sample recoveries and a detection limit of  $5 \times 10^{-6}$  g of creosote per m<sup>3</sup> of sample were achieved. HPLC. with either fluorescence or UV detection, is an EPA-recommended method for the analysis of both solid and liquid hazardous waste (EPA, 1986c). At present, HPLC does not achieve the high detection capability of capillary gas chromatography. HPLC however, does offer some advantages for the determination of creosote-derived PAH complex in environmental samples. Firstly, HPLC offers a variety of stationary phases capable of providing . unique selectivity for the separation of PAH isomers that are often difficult

to separate by GC. Secondly, ultraviolet absorption and fluorescence detection provide sensitive and more importantly, selective detection of PAHs in HPLC technique.

## 6.3 ADEQUACY OF THE DATABASE

Section 104(i)5 of CERCLA directs the Administrator of ATSDR (in consultation with the Administrator of EPA and agencies and programs of the Public Health Service) to assess whether adequate information on the health effects of creosote is available. Where adequate information is not available, ATSDR, in cooperation with the National Toxicology Program (NTP), is required to assure the initiation of a program of research designed to determine the health effects (and techniques for developing methods to determine such health effects) of creosote.

The following categories of possible data needs have been identified by a joint team of scientists from ATSDR, NTP, and EPA. They are defined as substance-specific informational needs that, if met would reduce or eliminate the uncertainties of human health assessment. In the future, the identified data needs will be evaluated and prioritized, and a substance-specific research agenda will be proposed.

#### 6.3.1 Identification of Data Needs

Methods for Determining Biomarkers of Exposure and Effect. Creosote is a complex mixture of organic compounds and no methods exist for measuring the parent compound in biological media. However, sensitive methods do exist for measuring components of the creosote mixture. Most of these methods involve detection of PAHs, the predominate components of creosote, and their metabolites. These analytical methods can reliably detect trace levels of PAHs in human tissues and body fluids, making them sensitive enough to measure background levels in the population, as well as levels at which biological effects might occur. PAHs, however, are not unique to creosote exposure. Development of analytical methods sensitive and selective enough to measure possibly unique or unusual components of creosote, or capable of yielding a unique "fingerprint" for the compound, would be useful in monitoring exposures that might occur in work environments and near hazardous waste sites where creosote has been detected.

The analytical methods for measuring PAHs and their metabolites in biological tissues and fluids are sensitive enough to measure levels at which health effects might occur, as well as background levels in the population. Methods also exist for measuring PAH-DNA adducts, and research efforts are underway to develop methods that will detect ultratrace levels of these adducts in biological media. The increased sensitivity may allow correlation between levels of these adducts and observed health effects of creosote-related PAH exposure. There is also a need for methods to quantitatively

correlate monitored levels of various PAHs in biological tissues or fluids to toxic effects in humans. Methods dependent on monitoring PAHs, however, are not specific for creosote exposure. Methods sensitive and selective enough to detect a more unique component or group of components making up the creosote mixture would allow a more accurate assessment of the health effects associated with exposure to monitored levels of creosote.

Methods for Determining Parent Compounds and Degradation Products in Environmental Media. Reliable and sensitive methods are available for measuring PAHs from creosote in soil, water, air, and other environmental media. Exposure to creosote is most likely to occur in industrial settings where creosote is manufactured or used. Creosote-contaminated water and soil are a concern in areas near hazardous waste sites and other areas where creosote might be concentrated. The analytical methods available are accurate and sensitive enough to quantitatively detect PAHs in these and other environmental media, and are effective for estimating creosote levels in media known to be contaminated with this compound. There is a lack of sensitive reliable methods for detecting and measuring creosote degradation products in environmental media. Development of such methods would allow assessment of the possible health effects of exposure to creosote metabolites and assist in determining the level of potential exposure to these products.

# 6.3.2 On-going Studies

No on-going studies concerning techniques for measuring and determining creosote in biological and environmental samples were reported.

# 7. REGULATIONS AND ADVISORIES

The international, national, and state regulations and advisories pertaining to creosote in air, water, and food are summarized in Table 7-1. Regulations and advisories pertaining to the components of creosote (i.e., PAHs and phenol) can be found in the ATSDR Toxicological Profiles for these chemicals (ATSDR 1989b, 1990). Creosote is on the list of chemicals appearing in "Toxic Chemicals Subject to Section 313 of the Emergency Planning and Community Right-to-Know Act of 1986" (EPA 1987b).

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# 7. REGULATIONS AND ADVISORIES

TABLE 7-1. Regulations and Guidelines Applicable to Creosote

gency	Description	Value	References
ARC	Carcinogenic Classification	Group 2A	IARC 1985
	<u>National</u>		
legulations			·
ir:			
SHA	Permissible Exposure Limit (PEL) Coal-tar Pitch	0.2 mg/m3	NIOSE 1985
	Volatiles (air)	•	
ater:	· ·		
PA OSW	Listing as hazardous waste from specific sources: bottom sediment sludge from the treatment of wastewaters from wood preserving processes that use creosote and/or pentachlorophenol	No data	EPA 1981b (40 CFR 261.33)
PA OERR	Reportable Quantity Notification required when discharges containing sludge are made into waterways	1 1b	EPA 1985 (50 FR 13474)
A	•		•
Other Media: CPA OPP	Classified for Restricted Use only by Certified Applicators, and for framing, piling applications and railroad tie repair only	No data	EPA 1986b (51 FR 1334)
	Listing as restricted use pesticide	No data	EPA 1986b (51 FR 1334)
PA OTS	Toxic chemical release reporting; community right-to-know	Yes	EPA 1987b
rsca	Prohibit application to wood intended for use in interiors or for use in contact with food, feed, or drinking water	No data	EPA 1984 (49 FR 28666)
Guidelines	100d, leed, of diliking wason	•	
ACGIH	Threshold Limit Value (TLV) 8 hour TWA Coal-tar Pitch Volatiles	0.2 mg/m3	ACGIH 1986
	(benzene soluble fraction)		
-	Human Carcinogen (coal-tar pitch volatiles)	Group Ala <sup>b</sup>	ACGIH 1986
NIOSH	10 hour TWA Coal-tar Pitch Volatiles (cyclohexane soluble fraction)	0.1 mg/m3	NIOSH 1985
	•		

Coal-tar Pitch Volatiles

# 7. REGULATIONS AND ADVISORIES

#### Table 7-1 (Continued)

Agency	Description	Value	References
State Regulations	and Guidelines		
State environmental agencies	Acceptable ambient air concentration guidelines and standards for several states		NATICH 1988
	Coal-tar Pitch Volatiles:		NATICH 1988
	Connecticut	2 μg/m3 (8 hr avg.)	•
	Kansas .	$0.0161 \mu g/m3$ (annual avg.)	
•	North Carolina	0 μg/m3 (24 hr avg.)	
	Nevada	0.005 μg/m3 (8 hr avg.)	
	Pennsylvania	0.48 μg/m3 (1 yr avg.)	
	Virginia	$2 \mu g/m3$ (24 hr avg.)	

<sup>a</sup>Group 2A -- Probable Human Carcinogen. <sup>b</sup>Group Ala -- Recognized Human Carcinogen.

ACGIH = American Conference of Governmental Industrial Hygiemists; EPA = Environmental Protection Agency; IARC = International Agency for Research on Cancer; IDLH = Immediately Dangerous to Life and Health; NATICH = National Air Toxics Information Clearinghouse; NIOSH = National Institute for Occupational Safety and Health; OERR = Office of Emergency and Remedial Response; OSHA = Occupational Safety and Health Administration; OPP = Office of Pesticide Programs; OSW = Office of Solid Waste; OTS = Office of Toxic Substances; PEL = Permissible Exposure Limit; TLV = Threshold Limit Value; TPQ = Threshold Planning Quantity; TWA = Time-Weighted Average; TSCA = Toxic Substances Control Act.

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#### 9. GLOSSARY

Acute Exposure -- Exposure to a chemical for a duration of 14 days or less, as specified in the toxicological profiles.

Adsorption Coefficient (Koc) -- The ratio of the amount of a chemical adsorbed per unit weight of organic carbon in the soil or sediment to the concentration of the chemical in solution at equilibrium.

Adsorption Ratio (Kd) -- The amount of a chemical adsorbed by a sediment or soil (i.e., the solid phase) divided by the amount of chemical in the solution phase, which is in equilibrium with the solid phase, at a fixed solid/solution ratio. It is generally expressed in micrograms of chemical sorbed per gram of soil or sediment.

Bioconcentration Factor (BCF) -- The quotient of the concentration of a chemical in aquatic organisms at a specific time or during a discrete time period of exposure divided by the concentration in the surrounding water at the same time or during the same period.

Cancer Effect Level (CEL) -- The lowest dose of chemical in a study, or group of studies, that produces significant increases in the incidence of cancer (or tumors) between the exposed population and its appropriate control.

Carcinogen -- A chemical capable of inducing cancer.

Ceiling Value -- A concentration of a substance that should not be exceeded, even instantaneously.

Chronic Exposure -- Exposure to a chemical for 365 days or more, as specified in the Toxicological Profiles.

Developmental Toxicity -- The occurrence of adverse effects on the developing organism that may result from exposure to a chemical prior to conception (either parent), during prenatal development, or postnatally to the time of sexual maturation. Adverse developmental effects may be detected at any point in the life span of the organism.

Embryotoxicity and Fetotoxicity -- Any toxic effect on the conceptus as a result of prenatal exposure to a chemical; the distinguishing feature between the two terms is the stage of development during which the insult occurred. The terms, as used here, include malformations and variations, altered growth, and in utero death.

EPA Health Advisory -- An estimate of acceptable drinking water levels for a chemical substance based on health effects information. A health advisory is

#### GLOSSARY

not a legally enforceable federal standard, but serves as technical guidance to assist federal, state, and local officials.

Immediately Dangerous to Life or Health (IDLH) -- The maximum environmental concentration of a contaminant from which one could escape within 30 min without any escape-impairing symptoms or irreversible health effects.

Intermediate Exposure -- Exposure to a chemical for a duration of 15-364 days as specified in the Toxicological Profiles.

Immunologic Toxicity -- The occurrence of adverse effects on the immune system that may result from exposure to environmental agents such as chemicals.

In Vitro -- Isolated from the living organism and artificially maintained, as in a test tube.

In Vivo -- Occurring within the living organism.

Lethal Concentration (LC<sub>LO)</sub> (LC<sub>LO)</sub> -- The lowest concentration of a chemical in air which has been reported to have caused death in humans or animals.

Lethal Concentration (50) (LC<sub>50</sub>) -- A calculated concentration of a chemical in air to which exposure for a specific length of time is expected to cause death in 50% of a defined experimental animal population.

Lethal  $Dose_{(LO)}$  (LD<sub>LO</sub>) -- The lowest dose of a chemical introduced by a route other than inhalation that is expected to have caused death in humans or animals.

Lethal  $Dose_{(50)}$  (LD<sub>50</sub>) -- The dose of a chemical which has been calculated to cause death in 50% of a defined experimental animal population.

Lethal  $Time_{(50)}$  (LT<sub>50</sub>) -- A calculated period of time within which a specific concentration of a chemical is expected to cause death in 50% of a defined experimental animal population.

Lowest-Observed-Adverse-Effect Level (LOAEL) -- The lowest dose of chemical in a study, or group of studies, that produces statistically or biologically significant increases in frequency or severity of adverse effects between the exposed population and its appropriate control.

Malformations -- Permanent structural changes that may adversely affect survival, development, or function.

Minimal Risk Level -- An estimate of daily human exposure to a chemical that is likely to be without an appreciable risk of deleterious effects (noncancerous) over a specified duration of exposure.

#### 9. GLOSSARY

Mutagen -- A substance that causes mutations. A mutation is a change in the genetic material in a body cell. Mutations can lead to birth defects, miscarriages, or cancer.

Neurotoxicity -- The occurrence of adverse effects on the nervous system following exposure to chemical.

No-Observed-Adverse-Effect Level (NOAEL) -- The dose of chemical at which there were no statistically or biologically significant increases in frequency or severity of adverse effects seen between the exposed population and its appropriate control. Effects may be produced at this dose, but they are not considered to be adverse.

Octanol-Water Partition Coefficient (Kow) -- The equilibrium ratio of the concentrations of a chemical in n-octanol and water, in dilute solution.

Permissible Exposure Limit (PEL) -- An allowable exposure level in workplace air averaged over an 8-hour shift.

 $q_1*$  -- The upper-bound estimate of the low-dose slope of the dose-response curve as determined by the multistage procedure. The  $q_1*$  can be used to calculate an estimate of carcinogenic potency, the incremental excess cancer risk per unit of exposure (usually  $\mu g/L$  for water, mg/kg/day for food, and  $\mu g/m^3$  for air).

Reference Dose (RfD) -- An estimate (with uncertainty spanning perhaps an order of magnitude) of the daily exposure of the human population to a potential hazard that is likely to be without risk of deleterious effects during a lifetime. The RfD is operationally derived from the NOAEL (from animal and human studies) by a consistent application of uncertainty factors that reflect various types of data used to estimate RfDs and an additional modifying factor, which is based on a professional judgment of the entire database on the chemical. The RfDs are not applicable to nonthreshold effects such as cancer.

Reportable Quantity (RQ) -- The quantity of a hazardous substance that is considered reportable under CERCLA. Reportable quantities are (1) 1 lb or greater or (2) for selected substances, an amount established by regulation either under CERCLA or under Sect. 311 of the Clean Water Act. Quantities are measured over a 24-hour period.

Reproductive Toxicity -- The occurrence of adverse effects on the reproductive system that may result from exposure to a chemical. The toxicity may be directed to the reproductive organs and/or the related endocrine system. The manifestation of such toxicity may be noted as alterations in sexual behavior, fertility, pregnancy outcomes, or modifications in other functions that are dependent on the integrity of this system.

#### 9. GLOSSARY

Short-Term Exposure Limit (STEL) -- The maximum concentration to which workers can be exposed for up to 15 min continually. No more than four excursions are allowed per day, and there must be at least 60 min between exposure periods. The daily TLV-TWA may not be exceeded.

Target Organ Toxicity -- This term covers a broad range of adverse effects on target organs or physiological systems (e.g., renal, cardiovascular) extending from those arising through a single limited exposure to those assumed over a lifetime of exposure to a chemical.

Teratogen -- A chemical that causes structural defects that affect the development of an organism.

Threshold Limit Value (TLV) -- A concentration of a substance to which most workers can be exposed without adverse effect. The TLV may be expressed as a TWA, as a STEL, or as a CL.

Time-Weighted Average (TWA) -- An allowable exposure concentration averaged over a normal 8-hour workday or 40-hour workweek.

Toxic Dose  $(TD_{50})$  -- A calculated dose of a chemical, introduced by a route other than inhalation, which is expected to cause a specific toxic effect in 50% of a defined experimental animal population.

Uncertainty Factor (UF) -- A factor used in operationally deriving the RfD from experimental data. UFs are intended to account for (1) the variation in sensitivity among the members of the human population, (2) the uncertainty in extrapolating animal data to the case of human, (3) the uncertainty in extrapolating from data obtained in a study that is of less than lifetime exposure, and (4) the uncertainty in using LOAEL data rather than NOAEL data. Usually each of these factors is set equal to 10.

#### APPENDIX

#### PEER REVIEW

A peer review panel was assembled for creosote. The panel consisted of the following members: Dr. David Strayer, Health Science Center, University of Texas; Dr. Paul Mushak, Private Consultant, Durham, NC; and Dr. Rick Irvin, TEES Engineering Toxicology Division, Texas A&M University. These experts collectively have knowledge of creosote's physical and chemical properties, toxicokinetics, key health end points, mechanisms of action, human and animal exposure, and quantification of risk to humans. All reviewers were selected in conformity with the conditions for peer review specified in Section 104(i)(13) of the Comprehensive Environmental Response, Compensation, and Liability Act, as amended.

A joint panel of scientists from ATSDR and EPA has reviewed the peer reviewers' comments and determined which comments will be included in the profile. A listing of the peer reviewers' comments not incorporated in the profile, with a brief explanation of the rationale for their exclusion, exists as part of the administrative record for this compound. A list of databases reviewed and a list of unpublished documents cited are also included in the administrative record.

The citation of the peer review panel should not be understood to imply its approval of the profile's final content. The responsibility for the content of this profile lies with the Agency for Toxic Substances and Disease Registry.

# U.S. ENVIRONMENTAL PROTECTION AGENCY REGION 4, SCIENCE and ECOSYSTEM SUPPORT DIVISION ATHENS, GA 30605-2720

4SESD-EIB

JUL 2 2 2003

Reference 17 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423

# **MEMORANDUM**

SUBJECT:

Gulf States Creosoting Sampling Event Report

SESD Project No. 03-0474, 03-0558

FROM:

Brian Striggow, Environmental Protection Specialist

Superfund and Air Section

THRU:

Don Hunter, Acting Chief

Superfund and Air Section

TO:

Brian Farrier

Waste Management Division South Site Management Branch

Attached please find the Sampling Event Report for the Gulf States Creosoting site in Flowood, Mississippi. If you have any questions about the subject document, please contact me at (706) 355-8619 or at email striggow.brian@epamail.epa.gov.

### Attachment

cc:

Shanieka Pennamon, Weston Solutions, Inc.

# **Gulf States Creosoting Flowood, Mississippi**

Sampling Event Report July 2003

United States Environmental Protection Agency Region 4 Science and Ecosystem Support Division Enforcement and Investigations Branch Superfund and Air Section

# GULF STATES CREOSOTING FLOWOOD, MISSISSIPPI

# Introduction

At the request of the US EPA Region 4, Waste Management Division (WD), South Site Management Branch (SSMB), the Science and Ecosystem Support Division (SESD) Enforcement and Investigations Branch (EIB) conducted a sampling event at the Gulf States Creosoting site in Flowood, Mississippi. The work was conducted the week of April 21, 2003

The purpose of the sampling event was to collect and analyze environmental samples for a Site Investigation (SI) report to be prepared by an EPA Superfund Technical Assistance Response Team (START) contractor. The information presented here is intended only to support the SI report preparation effort - it is not a comprehensive evaluation of the site.

The following materials are attached.

Title	Description		
Appendix A	Figures		
Figure 1	Sampling Locations		
Figure 2	Background Locations		
Appendix B Tables			
Table 1	Sample Descriptions and Coordinates		
Table 2	Surface Soil Analytical Results, positive values only		
Table 3	Subsurface Soil Analytical Results, positive values only		
Table 4	Sediment Analytical Results, positive values only		
Table 5	Groundwater Analytical Results, positive values only		
Table 6	Surface Soil Analytical Results		
Table 7	Subsurface Soil Analytical Results		
Table 7	Sediment Analytical Results		
Table 9	Groundwater Analytical Results		
Table 10	Surface Soil Miscellaneous Analytical Results		
Table 11	Subsurface Soil Miscellaneous Analytical Results		
Appendix C	Site Photographs		
Appendix D	Log Book Photocopies		
Appendix E Laboratory Data Sheets			

# **Background**

The Gulf States Creosoting Site site is located in Flowood, Mississippi, approximately 2 miles East of Jackson, Mississippi. The creosoting site operations ceased around 1950 and the site is currently in agricultural and industrial use. A levee separates the west side of the site from several oxbow lakes of the Pearl River identified on USGS maps as 'Creosote Slough'. The present levee was constructed during or after the creosoting operations. There is no visible evidence of the previous creosoting operations at or near the site.

# Sampling and Methodology Summary

All samples were collected and handled in accordance with the US EPA, Region 4, SESD, Environmental Investigation Standard Operating Procedures and Quality Assurance Manual, Nov, 2001 (EISOPQAM). Samples were analyzed in accordance with the Laboratory Operations and Quality Control Manual, Jan. 2002, and/or the Contract Laboratory Program (CLP) Statement of Work.

Sampling locations were designated by the Start Contractor, Weston Solutions, Inc., after a joint site reconnaissance with EPA, Weston and State of Mississippi personnel. The site and surrounding area, with sample locations indicated, is shown in Figure 1. The control sample locations are shown in Figure 2. Control samples of all media (GS-01-GW,SS,SB,SD) were collected northeast of the site behind Jackson Preparatory School on Hwy. 25. Additional control sediment samples were collected upstream of the site on the Pearl River (GS-06-SD) and on Prairie Branch (GS-08-SD), to isolate various potential influences.

All samples were analyzed for Volatile Organic Compounds(VOCs), Extractable (Semi-volatile) Organic Compounds (SVOCs), pesticides/PCBs, and total metal. The sediment samples were also analyzed for Total Organic Carbon (TOC) to give an indication of the level of natural organic matter. The sediments had an additional low-level extractables scan performed that is generally more suitable for the high-moisture content sediments.

All soil samples were collected with a hand auger and mixed in glass pans after collection of VOC samples in Encore samplers. Surface soils (designated with a SS suffix) were collected from a 0 to 6" Below Land Surface (BLS) sampling interval. With the exception of two samples, all subsurface soils (designated with a SB suffix) were collected from a 24"-36"BLS interval. The GS-10-SB and GS-14-SB samples were collected from a 12"-24" BLS interval. Note that the sampling plan called for GS-13-SB to be sampled from the shorter interval, but the GS-14-SB location was inadvertently sampled to the shallower depth instead. A co-located duplicate of sample GS-20-SS was collected and recorded as sample GS-20-SD. The analytical results for the duplicate are reported with results for other surface soil samples.

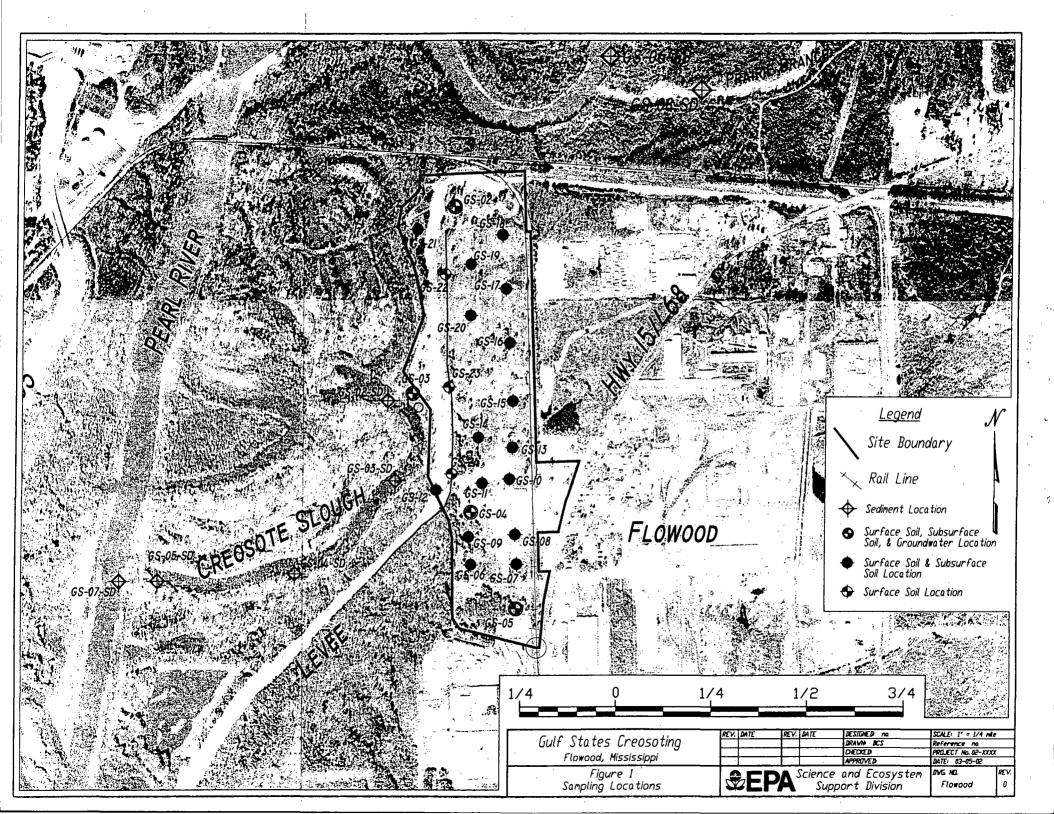
The sediment samples (designated with a SD suffix, except for duplicate soil sample GS-20-SD) were collected with a stainless-steel scoops and 10 foot handle-extension by samplers wading the sloughs or streams. The scoop was used to retrieve samples of the upper sediments (estimated 0-2") in the areas of interest. After collection, sediment samples were subsequently processed similarly to the soil samples.

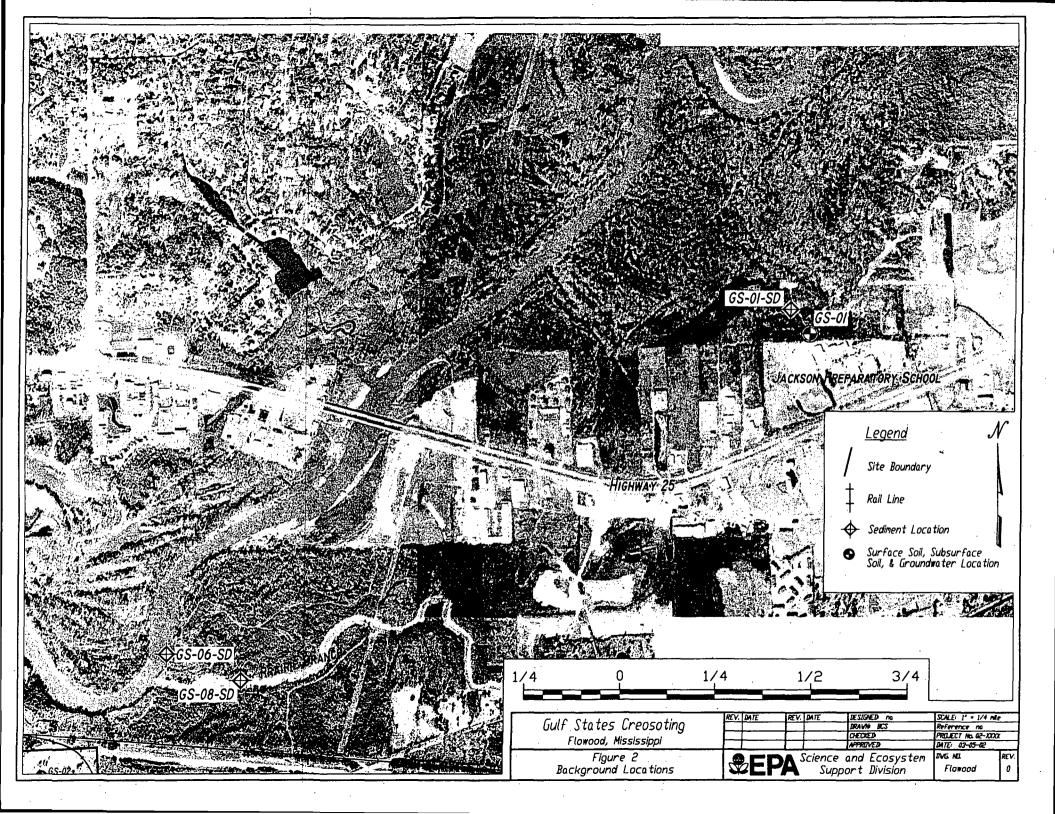
Temporary wells for goundwater sampling were installed using direct-push techniques. Geoprobe Systems® SP15 direct-push well screens were installed in the 16'-20' BLS interval at each location. Using low-flow techniques, each well was purged to a turbidity near or below 10 Nephelometric Turbidity Units (NTUs) prior to sampling.

At one location near Creosote Slough, creosote odors were detected on the recon/planning trip. This area was sampled as GS-03-SS,SB,and GW in this sampling effort. Creosote odors were not noted during sampling at this location or any other soil or groundwater sampling location. However, at sediment locations GS-02-SD and GS-03-SD a sheen was noted on the sample material in the pan and creosote odors were noted at location GS-03-SD. Note that while GS-XX-SS, SB, and GW are co-located, the GS-XX-SD samples are a separately numbered series at different locations from the soil and groundwater samples.

A preservative blank was processed after groundwater sampling was completed. Trip blanks were also shipped with each VOC sample shipment. No analytes or compounds were detected in any of these QA samples.

Appendix A Figures





# Appendix B Tables

**Table 1. Sample Descriptions and Coordinates** 

# Soils

		North	West
Sample	Description	Latitude	Longitude
GS-01-SS,SB	Background surface (0-6") and subsurface (24-36") soils from Jackson Prep	32.334569	90.109869
GS-02-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.318634	90.144468
GS-03-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.313000	90.146103
GS-04-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.309347	90.144088
GS-05-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.306379	90.142552
GS-06-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.307751	90.144129
GS-07-SS,SB	Surface (0-6") and subsurface (24-36") soils	r 32.307740	90.142494
GS-08-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.308640	90.142527
GS-09-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.308601	90.144193
GS-10-SS,SB	Surface (0-6") and subsurface (6-12") soils	32.310344	90.142697
GS-11-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.310225	90.143673
GS-12-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.310056	90.145333
GS-13-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.311298	90.142564
GS-14-SS,SB	Surface (0-6") and subsurface (6-12") soils	32.311612	90.143780
GS-15-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.312698	90.142520
GS-16-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.314473	90.142586
GS-17-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.316131	90.142686
GS-18-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.317754	90.142774
GS-19-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.316894	90.143933
GS-20-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.315338	90.143966
GS-20-SD	Co-located duplicate of GS-20-SS	32.315338	90.143966
GS-21-SS,SB	Surface (0-6") and subsurface (24-36") soils	32.317983	90.145814
GS-22-SS	Surface soils(0-6*) from drainage ditch	32.316609	90.144854
GS-23-SS	Surface soils(0-6") from drainage ditch	32.313169	90.144798
GS-24-SS	Surface soils(0-6") from drainage ditch	32.309375	90.145108

# Sediments

	North	West
Description	Latitude	Longitude
Background sediment from slough @ Jackson Prep School	32.335514	90.110716
Bottom sediment from Creosote Slough	32.312798	90.146997
Bottom sediment from Creosote Slough	32.310375	90.146789
Bottom sediment from Creosote Slough	32.307611	90.150453
Bottom sediement from Creosote Slough @ Pearl River	32.307500	90.155336
Background sediment from Pearl River	32.323095	90.138835
Bottom sediment from Pearl River	32.307525	90.156702
Background sediment from Prairie Branch	32.322002	90.135580
	Background sediment from slough @ Jackson Prep School Bottom sediment from Creosote Slough Bottom sediment from Creosote Slough Bottom sediment from Creosote Slough Bottom sediment from Creosote Slough @ Pearl River Background sediment from Pearl River Bottom sediment from Pearl River	DescriptionLatitudeBackground sediment from slough @ Jackson Prep School32.335514Bottom sediment from Creosote Slough32.312798Bottom sediment from Creosote Slough32.310375Bottom sediment from Creosote Slough32.307611Bottom sediment from Creosote Slough @ Pearl River32.307500Background sediment from Pearl River32.323095Bottom sediment from Pearl River32.307525

# Groundwater

		North	West
Sample	Description	Latitude	Longitude
GS-01-GW	Background sample at Jackson Prep School	32.334569	90.109869
GS-02-GW	Field Sample	32.318634	90.144468
GS-03-GW	Field Sample near Creosote Slough	32.313006	90.146002
GS-04-GW	Field Sample	32.309347	90.144088
GS-05-GW	Field Sample	32.306379	90.142552
GS-06-GW	Field Sample	32.307751	90.144129

Table 2. Surface Soil Analytical Results, positive values only. Page 1 of 6

	Units	GS01SS		GS02SS		GS03SS		GS04SS	Γ	GS05SS		GS06SS		GS07SS		GS08SS		GS09SS	П
% Moisture	%	15		23		25		23		15		26		17		5		19	
Aluminum	MG/KG	4700	J	7800	J	8900	J	4300	J	3100 J	J	6700	J	2800	J	1800	J	2500	J
Antimony	MG/KG		UJ	0.64	R	1	U		U	L	J		υT		c	1	U	1.2	R
Arsenic	MG/KG	6.3		3.3	J	11		2		13		2		3.6		3.4		0.95	R
Barium	MG/KG	72		120		67		60		37		85		34		21	1	25	
Beryllium	MG/KG	0.33		0.94		0.94		0.37		0.3		0.28		0.2		0.19		0.15	
Cadmium	MG/KG		Ü		U		U		U	L	7		U		U	0.09			U
Calcium	MG/KG	1100		690		790		690		500	$\neg$	1200		810		800		340	
Chromium	MG/KG	8.4		. 6.1		20		6		7.5	$\top$	7.7		4.5		4		4.1	
Cobalt	MG/KG	5.8		17		8.1		1.7		6.6	$\neg \top$	1.4		2		1.5		0.95	
Copper	MG/KG	6		2	J		U		C	l l	) I		ŪΠ		Ü		U	•	IJ
tron	MG/KG	10000		8400		27000	J	4500	J	15000 J	7	5600	J	5400	J	4700	J	2800	
Lead	MG/KG	11	_	17		120		18		19	$\neg$	17		11		12		8.8	
Magnesium	MG/KG	520	ے	470	J	680		. 280		180	$\neg$	550		250		180		160	J
Manganese	MG/KG	630		2800		550		140		610		170		180		230		28	
Potassium	MG/KG	380	J	240	J	350		200		120		430		170		150		120	J
Selenium	MG/KG	1.3		0.93	R	1.9	J		٥	Ü	J		ŪŢ		U	1	U		U
Silver	MG/KG	0.39	R	0.71	R	0.63		0.26	R	0.43		0.29			U		U	0.28	R
Sodium	MG/KG	270		300		82	J	120	J	58 J	厂	110		63			U	300	
Thallium	MG/KG		U		U		U		U	u	7	1	Ū		U		U		U
Vanadium	MG/KG	17		15		46		11		18		13		8.1		7.1		6.6	
Zinc	MG/KG	29		47		34		27		23		35		24		26		19	

	Units	GS01SS	GS02SS	GS03SS	GS04SS	GS05SS	GS06SS	GS07SS	GS08SS	GS09SS
% Moisture	%	14	23	. 24	24	15	26	15	4	26
4,4'-DDE (p,p'-DDE)	UG/KG	U	U	U	Û	U	U	U	U	U
4,4'-DDT (p,p'-DDT)	UG/KG	U	U	U	U	U	U	12 J	41 J	Ų
alpha-BHC	UG/KG	U	U	U	U	U	U	U	U	U
Endrin	UG/KG	U	U	U	· U	U	Ü	U	U	U
Endrin Aldehyde	UG/KG	Ü.	U	Ü	U	18 J	U	· U	17 J	U
Endrin Ketone	UG/KG	U.	U	U	U	U	U	U	U	U
Methoxychlor	UG/KG	U	U	U	U	U	U	U	74 NJ	· U

# Extractables Scan Page 2 of 6

	Units	GS01SS	GS02SS	GS03SS	GS04SS	GS05SS	GS06SS	GS07SS	GS08SS	GS09SS
% Moisture	%	14	23	24	. 24	15	26	15	4	26
1,1-Biphenyl	UG/KG	U	U	U	U	L	J U	45 J	U	U
2-Methylnaphthalene	UG/KG	U	U	U	0	l l		120 J	U	U
Acenaphthene	UG/KG	U	. U	U	J	(		170 J	U	U.
Acenaphthylene	UG/KG	U	U	U	٥	120 5	U	320 J	480 J	U
Anthracene	UG/KG	Ü	U	U	U	470	U	970	930 J	U
Benzo(a)Anthracene	UG/KG	U	U_	U	J	770	U	2500	1500	U
Benzo(b)Fluoranthene	UG/KG	U	U	U	ט	1800 J		4600 J	11000 J	U
Benzo(ghi)Perylene	UG/KG	U	U	U	J	380 J		690	3800	U
Benzo(k)Fluoranthene	UG/KG	· U	U	U	Ú	2100 J	U	5200 J	12000 J	U
Benzo-a-Pyrene	UG/KG	U	U	. U	U	810	U	2100	3700	U
Carbazole	UG/KG	U	U	U	U	. 200 J	U	510	190 J	U
Chrysene	UG/KG	[U]	U	U	U	1500	U	4900	2300	U
Dibenzo(a,h)Anthracene	UG/KG	UJ	UJ	Ų	J	190 J	U	520	1700	U
Dibenzofuran	UG/KG	U	· U	U	Ü		U	240 J	U	U
Fluoranthene	UG/KG	U	U	Ú	U	2600	U	8200	2600	U
Fluorene	UG/KG	U	U	U	U	l l	U	180 J	U	
Indeno (1,2,3-cd) Pyrene	UG/KG	UJ	UJ	U	U	610	U	1100	5000	U
Naphthalene	UG/KG	U	U	Ų	· U	į.	U	87 J	U	U
Pentachiorophenoi	UG/KG	UJ	UJ	U	U	1	U	U	U	U
Phenanthrene	UG/KG	U	U	Ú	U	320 J	U	4200	610 J	U
Pyrene	UG/KG	. U	U	Ū	U	1500	U	6200	2500	U

# Volatiles Scan

	Units	GS01SS		GS02SS	Π	GS03SS		GS04SS	GS05SS		GS06SS	GS07SS		GS08SS	Ť	GS09SS	
% Moisture	%	14		23		24		24	-15		. 26	15		1	NR	26	
Acetone	UG/KG	100	J	180	J	25	J	28 J	91	J	82 J	87	J		NR		וַט
Methyl Acetate	UG/KG		U		U		Ü	U		Ų	U		U		NR		U
Methyl Ethyl Ketone	UG/KG		C	21	l _		Ü	UJ		ŪĴ	UJ		UJ		NR		IJ

Table 2. Surface Soil Analytical Results, positive values only. Page 3 of 6

	Unite	GS10SS	GS11SS	GS12SS		GS13SS	GS14SS	1	GS15SS	GS16SS	GS1	755		GS18SS	
% Moisture	%	11	7	18		11	. 14		15	21		21	-:-	22	
Aluminum	MG/KG	4700 J	4400 J	5500	J	4400 J	4800	J	8000 J	7700 J	9	000	J	4900	J
Antimony	MG/KG	1 R	W		U	0.6 R	0.69	J	UJ	T.	JJ		UJ		UJ
Arsenic	MG/KG	6.8	3.7	4.4		6.3	9		6.9	5.5	1	5.5		5.4	
Barium	MG/KG	120	54	83		58	65		72	50		120		53	
Beryllium	MG/KG	0.46	0.44	0.46		0.54	0.44		0.71	0.43		0.94		0.35	
Cadmium	MG/KG	0.34	0.19		Ü	0.16 R	0.11	R	0.1		j i	0.07	R		U
Calcium	MG/KG	2600	660	760		1200	1800		2000	1200	2	300		890	
Chromium	MG/KG	23	12	8.9		7.8	8.6		11	10		8		9.9	
Cobalt	MG/KG	4.8	5.4	7.3		4.2	7.3		11	4		12		7.6	
Copper	MG/KG	19 J	UJ		υ	7.2 J		UJ	ÜJ		J	$\neg \neg$	U		Ū
Iron	MG/KG	14000	11000	11000	J	12000	11000		12000	12000	12	000		10000	$\Gamma$
Lead	MG/KG	80	290	25		33	24		26	15 J		20	J	15	J
Magnesium	MG/KG	780 J	260 J	380		280 J	340	J	430 J	510 J		490	J	340	J
Manganese	MG/KG	500	880	990		940	870	· .	1800	240	2	500		590	
Potassium	MG/KG	560 J	330 J	280		390 J	200	J	580 J	490 J		620	J	180	J
Selenium	MG/KG	0.98 R	0.77 R		U	1.2	0.63	R	1.1 R	1.3		1.5	J	1.1	
Silver	MG/KG	0.49	0.38	0.43		0.33 R	0.37		0.5	0.39 F	3 (	0.51	R	0.29	
Sodium	MG/KG	320	200	91		230	300		310	270		280		310	
Thallium	MG/KG	U	U		Ū	U		U ,	U	l	j		υ		U
Vanadium	MG/KG	18	16	22		17	21		23	22		23		20	
Zinc	. MG/KG	160	840	28		72	44		48	33		65		34	

	Units	GS10SS	GS11SS	GS12SS	GS13SS	GS14SS		GS15SS	GS16SS	GS17\$S	GS18SS
% Moisture	%	12	7	17	13	14	$\neg$	16	21	21	23
4,4'-DDE (p,p'-DDE)	UG/KG	U	U	U	U		Ú	U	2.5 J	U	Ü
4,4'-DDT (p,p'-DDT)	UG/KG	43 J	30 J	U	38 J	10	NJ	30 NJ	Ü	U	Ū
alpha-BHC	UG/KG	U	U	C	U		Ū	U	U	U	U
Endrin	UG/KG	U	Ú	U	42 J		υ	U	U	6.4 J	U
Endrin Aldehyde	UG/KG	U	U	U	49 J		U	27 NJ	Ü	U	Ü
Endrin Ketone	UG/KG	U	Ú	U	150		U ·	U	14 J	20 J	Ü
Methoxychlor	. UG/KG	81 NJ	U	· U	U	34	NJ	U	U	U	U

# Extractables Scan Page 4 of 6

. 490 4 01 0	Units	GS10SS	$\neg$	GS11SS	GS12SS	GS13SS	~	GS14SS		GS15SS		GS16SS	$\Box$	GS17SS		GS18SS	Ŀ
% Moisture	%	12		7	17	13	$\neg$	14		16		21		21		23	
1,1-Biphenyl	UG/KG	- 1	Ü	J	Ú		J.		U		Ū		Ū		U		Ū
2-Methylnaphthalene	UG/KG	69	J	70 J	U	L	7		υ		Ū		U		U		U
Acenaphthene	UG/KG	97 .	J	U	U	Ĺ	_		U		Ū		J		J		Ū
Acenaphthylene	UG/KG	1100	_ 7	330 J	U	1000 J			٥	250	J	87	J	270	J		U
Anthracene	UG/KG	1700		830	U	1600		1100		370	J		U	380	J		U
Benzo(a)Anthracene	UG/KG	2800	.	1000	U	4300		2300		690		260	J	1000			U
Benzo(b)Fluoranthene	UG/KG	5200		2500	U	7300 J	ī —	4200		1500		700	J	2600	7		Ū
Benzo(ghi)Perylene	UG/KG	1300		460	U	2000		820		360	J	92	J	.290	J		Ū
Benzo(k)Fluoranthene	UG/KG	2300		1700	Ü	5400 J		. 2400		1100		730	j	2700	J		Ū
Benzo-a-Pyrene	UG/KG	3100		800	U	3100		2300		650		230	J	1000			Ü
Carbazole	UG/KG	480		210 J	· U	540 J		280	J	250	J	62	J	130	Ĵ_		Ū
Chrysene	UG/KG	4300	7	1500	U	6200		3100		1700	_	390	J	1600			Ü
Dibenzo(a,h)Anthracene	UG/KG	750	7	310 J	U	1100		580		200	J	95	J	330	J		U
Dibenzofuran	UG/KG	120	ĵΠ	140 J	U	L	, ,		U		Ū		Ü		U		Ū
Fluoranthene	UG/KG	4700		1900	U	11000		3100		3100		760		2500			Ü
Fluorene	UG/KG		ũΠ	U	U	L	,		U		Ū		Ü		U		U
Indeno (1,2,3-cd) Pyrene	UG/KG	2300	$\neg$	940	U	3100	П	1700		740	J	270	J	890			U
Naphthalene .	UG/KG	110	)	120 J	U	i.	П		U		Ū		U		U		U
Pentachlorophenol	UG/KG	690	Г	640 J	U	- L	П		U		ŪĴ		U		U		U
Phenanthrene	· UG/KG	1400		540	U	1600		320	J	1100		310	J	250	J		Ū
Pyrene	UG/KG	5000		1700	U	6800		3300		2200		540		1700			Ū

# Volatiles Scan

				<u>.</u>												
	Units	GS10SS		GS11SS	GS12SS	GS13SS		GS14SS		GS15SS	T	GS16SS		GS17SS		GS18SS
% Moisture	%	12		7	17	13		14		16	丁	21		21		23
Acetone	UG/KG	84	j	220 J	UJ	200 J		75	7	100 J		56	J	260 J	ĵ ,	130 J
Methyl Acetate	UG/KG		Ū	U	U	l	1		Ü	3 J	7	1	J	1	U	U
Methyl Ethyl Ketone	UG/KG		J	21 J	14 J	15 J			U	12	7		J	28 J	,	14 J

Table 2. Surface Soil Analytical Results, positive values only. Page 5 of 6

	Units	GS19SS	_	GS20SS		GS20SD		GS21SS		GS22SS		GS23SS		GS24SS	
% Moisture	%	15		20		19		22		- 14		26		47	
Aluminum	MG/KG	8500	5	6700	J	6100	J	6800	J	4200	J	7500	5	4900	J
Antimony	MG/KG		ÚĴ	· i	UJ	0.59	IJ		UJ		c		IJ	_	U
Arsenic	MG/KG	6.5		5.8		4.5		2.7		4		5		6.5	
Barium	MG/KG	89		410		130		88	1	42		83		120	
Beryillum	MG/KG	0.73		1.2		0.76		0.81		0.43		0.5		0.61	
Cadmlum	MG/KG		U	0.17		0.05	Ú		U.		5	0.1	R	0.16	
Calcium	MG/KG	2000		1700		1400		190		280		960	•	1600	
Chromium	MG/KG	8.3		8.2		7.2		5.6		9.4		12		12	
Cobalt	MG/KG	13		: 17		6.8		14		5.5		6.8		6.8	
Copper	MG/KG		U		Ū	4.3	U		U		Ų	12	J		U
Iron	MG/KG	12000		12000		8800		7300		9700	J	13000		17000	J.
Lead	MG/KG	19	J.	24	J	. 22	J	13	J	9.8		24		28	
Magnesium	MG/KG	530	J	380	J	410	J	270	J	220		630	J	450	
Manganese	MG/KG	1900		5500		1800		2200		470		920		890	
Potassium	MG/KG	550	J	240	J	250	J	250	J	170		480	J	370	
Selenium	MG/KG	1.2		1.3	L.	1.1	R	0.8			Ū	1.1	R		U
Silver	MG/KG	0.53	R	0.9	R	0.42		0.52		0.2		0.46	R	0.47	
Sodium	MG/KG	290		290		290	-	300			Ú	340		86	J
Thailium	MG/KG		Ü	1.5		0.77	U		Ü		Ū		Ū		Ū
Vanadium	MG/KG	22		22		17		13		21		23		26	$\overline{}$
Zinc	MG/KG	52		88		63		24		19		49		89	

	Units	GS19SS		GS20SS		GS20SD		GS21SS	$\Box$	GS22SS		GS23SS		GS24SS	Г
% Moisture	%	17		21		. 20		22		13		26		45	
4,4'-DDE (p,p'-DDE)	UG/KG		U	4.1	NJ	3.3	NJ		Û		U		U_		U
4,4'-DDT (p,p'-DDT)	UG/KG		υŢ		U		U		Ū		Ū		U		U
alpha-BHC	UG/KG		U		J		U		U		U	1.7	NJ		U
Endrin	UG/KG	· · · · · ·	ŪŢ		Ü		U		Ū		U		U		U
Endrin Aldehyde	UG/KG		U	1	U		U		U		U		U		U
Endrin Ketone	UG/KG	~	U		U		Ü		U		U	6.9	J_		U
Methoxychlor	UG/KG		U		U		U		ŪΪ		U		U		Ū

# Extractables Scan

	Units	GS19SS		GS20SS		GS20SD		GS21SS		GS22SS	G\$23\$\$	_	GS24SS	:T
% Moisture	%	17		21		20	_	22		13	26		45	T
1,1-Biphenyl	UG/KG		U		U	T	Ū		U	U		U		Ū
2-Methylnaphthalene	UG/KG		Ü.		U	1	Ü		Ū	U		Ü		U
Acenaphthene	UG/KG		U		Ü		Ü		Ū	U		U		U
Acenaphthylene	UG/KG		U		J		U		Ū	U	66	J	150	J
Anthracene	UG/KG		U	1	U		c		Ū	U		U	330	J
Benzo(a)Anthracene	UG/KG		U	58	3	44	J		Ū	Ü	430	J	990	$\top$
Benzo(b)Fluoranthene	UG/KG		Ü	160	J	140	5		Ū	Ū		NA	2700	J
Benzo(ghi)Perylene	UG/KG		U		U				Ū	U	180	J	330	J
Benzo(k)Fluoranthene	UG/KG		U	170	5	150	J		Ũ	U	520		2600	15
Benzo-a-Pyrene	UG/KG		υ	. 46	J	47	J		Ū	U	380	J	870	Т
Carbazole	UG/KG		U ,		٦				Ū	U		U		U
Chrysene	UG/KG		U_	98	5	69	J		Ŭ	U	510		1100	T
Dibenzo(a,h)Anthracene	UG/KG		U		5				Ū	U	120	J	260	J
Dibenzofuran	UG/KG		U		U				Ū	U		U		TU
Fluoranthene	UG/KG		U	73	J	81	5	7	Ū	U	520		1300	1
Fluorene	UG/KG		Ū		Ü				Ū	U		U		U
Indeno (1,2,3-cd) Pyrene	UG/KG		Ü	63	J	51	5		Ū	U	340	J	550	J
Naphthalene	UG/KG		U	1	Ü				Ū	U		U		U
Pentachlorophenol	UG/KG		U		Ū				U.	U		UJ		U
Phenanthrene	UG/KG		Ū		ΰ				Ū	U	59	J		U
Pyrene	UG/KG		Ū	100	J	63	J		U	U	450		1600	$\overline{}$

#### Volatiles Scan

·	Units	GS195S		Ι	GS20SS		GS20SD		GS2155		GS22SS		GS23SS		GS24SS	1
% Molsture	%	17	_		21		20		22		13		26		45	
Acetone	UG/KG	180	Ĵ	Ţ.	83	5	230	J	270	J	110	J	130	J	220	J
Methyl Acetate	UG/KG	3	J			υ		U		U		U		U		U
Methyl Ethyl Ketone	UG/KG	17	J	Ι.		U	23	J	31	J		UJ	14			UJ

### **Data Qualifiers**

U-Analyte not detected at or above reporting limit.

J-Identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolities of technical chlordane

Table 3. Subsurface Soil Analytical Results, positive values only. Page 1 of 6

	Units	GS01SB		GS02SB		GS03SB		GS04SB		GS05SB		GS06SB	GS07SB		GS08SB		GS09SB	
% Moisture	%	.18		19		28		27		23		27	26		18		20	
Aluminum	MG/KG	7300	J	7800	J	8600	J	6600	J	5300	J	7700 J	9400	J	8100	J.	5800	J
Antimony	MG/KG		UJ		UJ		U		Ū		U	U		U		Ü		ŲJ
Arsenic	MG/KG	6.7		3.4	J	9.4		1.9			U	2.7	2.7		3.5			Ū-
Barium	MG/KG	62		34		63	_	57		. 59		48	50		45		29	
Beryillum	MG/KG	0.31		0.36		0.64		0.41		0.27		0.24	0.35		0.59		0.35	
Cadmium	MG/KG		U		U		U		U		U	U		U		Ū		U
Calcium	MG/KG	850		170		1400		670		580		990	670		450		540	
Chromium	MG/KG	12		9.5		14	$\overline{}$	6.6		4.9	_	7.5	8		9.2		5.3	
Cobalt	MG/KG	3.3		2.6		4.3		1.4		0.78	R	1.2	1.6		9.4		1	R
Copper	MG/KG	6.4			5		U		v		Ū	U		U		U		IJ
Iron	MG/KG	13000		13000		22000	J.	7500	,	2300	J	5800 J	8200	J	12000	J	3800	
Lead	MG/KG	8.2	7	6.6		15		9.2		6.7		11	11		5.7	$\neg$	8.4	
Magnesium	MG/KG	580	J	480	J	770		370		220		580	550		370		290	J
Manganese	MG/KG	180		200		200		60		16	_	18	19		1000		9.3	
Nickel	MG/KG	- :	5		Ü		U	l	U		Ū,	U		Ū		U		Ü
Potassium	MG/KG	320	7	320	J	360		190		170		240	260		350		190	J
Selenium	MG/KG	0.7		1.4		1.6			Ū		U	U		Ū		U		U
Silver	MG/KG	0.38		0.38	R	0.49	R.		U		U	0.28 R		Ū	0.32			ΰ
Sodium	MG/KG	290		290		98	J	300		190	J	250	220		46		440	
Vanadium	MG/KG	22		20		33		14		6.7		17	18		22	$\neg$	8.4	
Zinc	MG/KG	18		20		23		15		7.3		12	15		25		12	

	Units	GS01SB	GS02SB	GS03SB	GS04SB	GS05SB	GS06SB	GS07SB	GS08SB	GS09SB
% Moisture	%	18	19	28	23	19	26	26	17	21
4,4'-DDE (p,p'-DDE)	UG/KG	U	U	U	U	- U	U	U	U	U
4,4'-DDT (p,p'-DDT)	UG/KG	U	U	U	U	U	U	U	U	· U
Endrin Ketone	UG/KG	U	U	U	U	U	U	U	U	U
gamma-BHC (Lindane)	UG/KG	U	U	U	U	U	U	U	U	U
PCB-1260 (Aroclor 1260)	UG/KG	U	U	U	U	U	U	U	U	U

# Extractables Scan Page 2 of 6

	Units	GS01SB		GS02SB	GS03SB		GS04SB	T	GS05SB	GS06SB		GS07SB		GS08SB	GS09SB
% Moisture	%	18		19	28		23	7	19	26	$\neg$	26		17	21
2-Methylnaphthalene	UG/KG	ì	U	U		7	U		U	U			U	U	U
Acenaphthene	UG/KG	}·	Ū	U	į	7	U	Т	U	U	П		U	U	U
Acenaphthylene	UG/KG		U	U	T.	,	U	T	U	U			C	U	U.
Anthracene	UG/KG		U	U	1	)	U		U	U			U	74 J	U
Benzo(a)Anthracene	UG/KG		U	U	l	J_	U	Ţ	U	U			Ū	120 J	U
Benzo(b)Fluoranthene	UG/KG		U	U			U	$\Box$	U	U		47	J	1000 J	U
Benzo(ghi)Perylene	UG/KG		U	U		J	U	Ι	U	U			C	320 J	U
Benzo(k)Fluoranthene	UG/KG		U	U	į	<i>J</i>	U	1	Ū	U		45	J_	980 J	U
Benzo-a-Pyrene	UG/KG		U	U		<u> </u>	U		U	U			U_	240 J	U
bis(2-Ethylhexyl) Phthalate	UG/KG		C	U	l	IJ	UJ	ıΤ	450 J	U	П	1	U	U	U
Carbazole	UG/KG		Ū	U	l	,	U	I	U	U			Ü	U	U
Chrysene	UG/KG		C	U	Ĺ	J_[	U	I	C	U			U	220 J	U
Dibenzo(a,h)Anthracene	UG/KG		U	UJ		)	U		U	U			Ü	150 J	ט
Dibenzofuran	UG/KG		U	U	l	,	U	I	U	U			U	U	U
Fluoranthene	UG/KG		U	U	L	J	U	I	U	U			U_	200 J	U
Indeno (1,2,3-cd) Pyrene	UG/KG		U	UJ		<i>_</i>	U	I	U	U			Ü	420	U
Naphthalene	UG/KG		U	U	l	,	U	Т	U	U		1	J	U	U
Pentachlorophenol	UG/KG		U	UJ	l	7	U	I	U	Ü	J		U	U	U
Phenanthrene	UG/KG		U	U	U	Į.	U	I	U	U			J	U	U
Pyrene	UG/KG		U	U	L L		U	$\perp$	Ú	Ü	$\Box$	l	J ]	230 J	U

# Volatiles Scan

•	Units	GS01SB		GS02SB	GS03SB		GS04SB		GS05SB		GS06SB		GS07SB		GS08SB		GS09SB	
% Moisture	%	18		19	28		23		19		26		26		17		21	
Acetone	UG/KG		5	12 J	35	J		UJ		UJ	21	J	U	IJ	99	J		UJ
Benzene	UG/KG		Ü	U		U		U		U		U	U	ГТ <u> </u>	- 1	U		U
Methyl Ethyl Ketone	UG/KG		IJ	U		บัง		UJ		屷		IJ	U	IJ	11 .	J		UJ

Table 3. Subsurface Soll Analytical Results, positive values only. Page 3 of 6

	Units	GS10SB	GS11SB		GS12SB	GS13SB	GS14SB		GS15SB		GS16SB		GS17SB		GS18SB	
% Moisture	%	18	18		17	15	15		.21		23		21		18	. [
Aluminum	MG/KG	8600	9600	J	3700 J	8400 J	4600	J	11000	J	12000	J	6700	J	7700	J
Antimony	MG/KG	0.96	٦ 1.6	R	U	0.58 R		UJ		UJ		UJ		UJ		UJ
Arsenic	MG/KG	13	6.6		4.6	4.7	7.4		8.1		7.3	$\Gamma$	1.5		4.1	
Barium	MG/KG	430	. 41		19	23	54		45	,	49		61		26	
Beryllium	MG/KG	1.3	0.49		0.3	0.34	0.48		0.59		0.49		0.39		0.31	
Cadmium	.MG/KG	0.87		U	\U_{\\	U		U		U		U		U		U
Calcium	MG/KG	9400	320		160	400	620	$\Gamma$	710		390		620		190	
Chromium	MG/KG	19	17		8.5	11	7.1		21		17		7		8.7	
Cobalt	MG/KG	12	3.2	R	1.4	1.7	12	T	2.9		2.1	T	1.5		3.4	
Copper	MG/KG	40	j.	UJ	U	U		UJ	7.	J	7.7			U		Û
Iron	MG/KG	20000	20000		13000 J	14000	14000		24000		23000		5900		12000	
Lead	MG/KG	71	6.7		6.7	5.1	27		9.2	-	8.7	J	9.1	J	8	J.
Magnesium	MG/KG	2400	680	J	170	470 J	250	J	640	J.	650	J	310	J	470	J
Manganese	MG/KG	1800	120		55	82	1000		410		52		57		76	
Nickel	MG/KG	20		U	U	U	1	U		U		U.		U		U
Potassium	MG/KG	970	430	J	170	410 J	180	J	490	J	470	J	280	J	220	J
Selenium	MG/KG	1.2	1.1	R	1.4	U		U	2.3	J	1.7	J		Ü		U
Silver	MG/KG.	0.76	1.1	R	0.47	0.39	0.42		0.65	-	0.45	1	0.32	R	0.32	R
Sodium	MG/KG	490	350		96	310	260	1	290		290	1	340		540	
Vanadium	MG/KG	30	33		26	22	23	1	36		33	$\Gamma$	13		18	
Zinc	MG/KG	290	44		8.4	20	20		36		25		12		16	

	Units	GS10SB	GS11SB	GS12SB	GS13SB	GS14SB	GS15SB	GS16SB	GS17SB	GS18SB
% Moisture	%	17	17	26	15	15	21	22	22	25
4,4'-DDE (p,p'-DDE)	UG/KG	U	U	U	6.9	U	U	U	Ü	U
4,4'-DDT (p,p'-DDT)	UG/KG	110 N	U	U	U	16 J	U	U	U	U
Endrin Ketone	UG/KG	180 N	U	Ü	U	U	U	Ü	U	U
gamma-BHC (Lindane)	UG/KG	Ü	U	U	U	U	U	1.3 J	U	1.7 J
PCB-1260 (Aroclor 1260)	UG/KG	U	U	U	U	U	U	70	U	42 J

# Extractables Scan Page 4 of 6

	Units	GS10\$B	GS*	1SB	GS12SB	G	S13SB		GS14SB		GS15SB	GS16SB	GS17SB	GS18SB
% Moisture	%	17		17	26		15		15		21	22	22	25
2-Methylnaphthalene	UG/KG	200	J	U		J		C		U	U	U	U	U
Acenaphthene	UG/KG	210	J	U	}	J		C		U	U	U	U	U
Acenaphthylene	UG/KG	6100		U		J .		٥	330	J	U	U	U.	U
Anthracene	UG/KG	12000		49 J		ر		C	690		U.	U	U	U
Benzo(a)Anthracene	UG/KG	23000		U		١	56		1400	·	U	U	U	Ū.
Benzo(b)Fluoranthene	UG/KG	37000		U		)	160	J	2400		U	U	41. J	U
Benzo(ghi)Perylene	UG/KG	5200		U	]	<u> </u>		U	540		U	U	U	Ū
Benzo(k)Fluoranthene	UG/KG	26000		46 J		) [	170	J	2000		U	U	43 J	Ū
Benzo-a-Pyrene	UG/KG	25000		U		J [		U	1400		U	U	U	U
bis(2-Ethylhexyl) Phthalate	UG/KG		U	U		]		U		U	U	U	U	UJ
Carbazole	UG/KG	1800		U		J		U	210	J	U	υ	U	U
Chrysene	UG/KG	35000		U		j T	91	J	2100		U	U	U	U
Dibenzo(a,h)Anthracene	UG/KG	3300		U		)		U	390		UJ	U	U	U
Dibenzofuran	UG/KG	340	J	Ū		]		U		U	U	U	U	U
Fluoranthene	UG/KG	28000		U		J	130	J	2200		U	U	U	Ū
Indeno (1,2,3-cd) Pyrene	UG/KG	20000		Ū		J .	51	J	1100		UJ	U	U	U
Naphthalene	UG/KG	390	J	U	1	J		Ū		U	U	U	U	U
Pentachlorophenol	UG/KG	680	J	U	1	J		U		U	UJ	U	U	U
Phenanthrene	UG/KG	2700		U		J		U	510		U	U	υ	U
Pyrene	UG/KG	37000		U			89	J	2400		U	U	U	U

# Volatiles Scan

	Units	GS10SB		GS11SB		GS12SB		GS13SB		GS14SB	GS15SB		GS16SB	GS17SB	GS18SB
% Moisture	%	17		17		26		15		15	21		22	22	25
Acetone	UG/KG	78	J	24	J		UJ		υJ	55 J	67	J	ÜJ	U	UJ
Benzene	UG/KG	2	J		C		U		U	U	1	Ū	U	U	U
Methyl Ethyl Ketone	UG/KG		ح		J		IJ		UJ	. U		U	UJ	U	UJ

Table 3. Subsurface Soil Analytical Results, positive values only. Page 5 of 6

	Units	GS19SB		GS20SB		GS21SB	
% Moisture	%	20		16		20	
Aluminum	MG/KG	12000	J	4700	J	8600	J
Antimony	MG/KG		IJ		IJ		S
Arsenic	MG/KG	8.3		2.3		4.5	
Barium	MG/KG	43		24		28	
Beryllium	MG/KG	0.46		0.3		0.3	
Cadmium	MG/KG		U	,	U		U
Calcium	MG/KG	190		120		140	
Chromium	MG/KG	12		4.7		14	
Cobalt	MG/KG	9.6		2.2		1.9	
Copper	MG/KG		U		C	3.6	
Iron	MG/KG	20000		6900		13000	
Lead	MG/KG	12	J	6	7	· 6.8	J
Magnesium	MG/KG	680	J	210	5	420	J
Manganese	MG/KG	600		140		95	
Nickel	MG/KG		U		U		U
Potassium	MG/KG	480	J	170	-	300	J_
Selenium	MG/KG	1.7	J		C	0.95	R
Silver	MG/KG	0.53		0.25	R	0.4	
Sodium	MG/KG	280		340		310	
Vanadium	MG/KG	31		14		21	
Zinc	MG/KG	: 30		11		16	

	Units	GS19SB		GS20SB		GS21SB	Г
% Moisture	%	21		18		20	$ extstyle  ag{7}$
4,4'-DDE (p,p'-DDE)	UG/KG	!	U		U		Ū
4,4'-DDT (p,p'-DDT)	UG/KG		Ü		U		U
Endrin Ketone	UG/KG	. !	U		U		U
gamma-BHC (Lindane)	UG/KG		Ü		U		U
PCB-1260 (Arocior 1260)	UG/KG	39	J		Ū		Ū

### Extractables Scan Page 6 of 6

•	Units	GS19SB	·	GS20SB		GS21SB	
% Moisture	%	21		18		20	
2-Methylnaphthalene	UG/KG	i	U		U		U
Acenaphthene	UG/KG		Ū		U		U
Acenaphthylene	UG/KG		U		U		U
Anthracene	UG/KG	!	U		U	·	כ
Benzo(a)Anthracene	UG/KG		Ú		U		٥
Benzo(b)Fluoranthene	UG/KG		U		U		5
Benzo(ghi)Perylene	UG/KG		υ	120	J_		ح
Benzo(k)Fluoranthene	UG/KG		Ü		U		ردا
Benzo-a-Pyrene	UG/KG		U		U		5
bis(2-Ethylhexyl) Phthalate	UG/KG		UJ		U		5
Carbazole	UG/KG		U		U		כ
Chrysene	UG/KG	i_	U		U		5
Dibenzo(a,h)Anthracene	UG/KG		U		U		υ
Dibenzofuran	UG/KG		U		Ū		Ų
Fluoranthene	UG/KG		U		U		Ų
Indeno (1,2,3-cd) Pyrene	UG/KG	1.	U		U		U
Naphthalene	UG/KG		U		Ū		U
Pentachlorophenol	UG/KG	- 1	U		Ú		Ū
Phenanthrene	UG/KG	1	U		Ų		U
Pyrene	UG/KG		U		Ū		Ū

#### Volatiles Scan

	Units	GS19SB		GS20SB	$\neg \neg$	GS21SB	
% Moisture	- %	21		18		20	_
Acetone	UG/KG	i	UJ		U		UJ
Benzene	UG/KG		U		U		Ū
Methyl Ethyl Ketone	UG/KG		UJ		U		UJ

#### **Data Qualifiers**

U-Analyte not detected at or above reporting limit.

J-identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-identification of analyte is acceptable; reported value may be blased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to interferences. | A-Analyte analyzed in replicate, Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Table 4. Sediment Analytical Results, positive values only. Page 1 of 3

·	Units	GS01SD	Γ	GS02SD	Γ	GS03SD		GS04SD		GS05SD		GS06SD	$\Box$	GS07SD		GS08SD	
% Moisture	%	61		56		84		. 77		54		20		22		27	
Aluminum	MG/KG	8200	J	4700	J	16000	J	9100	J	4900	J	450	J	140	J	510	J
Antimony	MG/KG		UJ		UJ		IJ		UJ	2.1	R	0.7	R		UJ		U
Arsenic	MG/KG	5.7		1.8	R	7.9	R	5.5		2.9	R		U		U		U
Barium	MG/KG	89		77		300		120		71		6.8		3.2		15	
Beryllium	MG/KG	0.39	<u> </u>	0.56		1.3		1.2	L	0.51			U.		U	0.1	
Calcium	MG/KG	1500		880		2000		2000	ì	1200		250			Ū	180	
Chromium	MG/KG	13		7.8		24		15		8.4			U		U	1.4	
Cobait	MG/KG	4.4		4.7		12		11		5.5		0.61	R	0.47		1	
Iron	MG/KG	15000		6400		24000		13000		8400		950		340		2200	J
Lead	MG/KG	12	J	20	J	41	J	21		. 12		1.8		1.2	J	3.6	
Magnesium	MG/KG	1000	J	410	J	1200	J	1000	ے	660	J	64	J		ŪJ	81	
Manganese	MG/KG	340		500		720		280		320		52		53		44	
Potassium	MG/KG	700	J	410	J	1100	J	811	J	480	J	90	J	52	J	56	
Selenium	MG/KG	1.9	R		U	4.2			Ŭ		Ü		U		U		U
Silver	MG/KG	0.68	R	0.45	R	1.1	R	1.1	R	0.71	R	0.24	R		U		U
Sodium	MG/KG	690		580		1500		1200		600		330		330		68	
Vanadium	MG/KG	24		13		41		23		13		1.6	R	0.57	R	2.4	
Zinc	MG/KG	36		33		130		84		39		3.6			U	4.6	

# Pesticides & Aroclors Scan

	Units	GS01SD	GS02SD		GS03SD	GS04SD	1	GS05SD	GS06SD	GS07SD	GS08SD
% Moisture	%	55	52		79	78		54	20	18	19
4,4'-DDE (p,p'-DDE)	UG/KG	U	7.9		U		U	U	U.	U	U
beta-BHC	UG/KG	U		Ū	12		U	U	U	U	U
Endosulfan II (beta)	UG/KG	U	6.5	J	U		Ū	U	U	U	U
Endosulfan Sulfate	UG/KG	U	10	N	14 NJ		U	U	U	U	. U

# Extractables Scan

	Units	GS01SD	GS02SD		GS03SD		GS04SD	Γ	GS05SD	GS06SD		GS07SD		GS08SD
% Moisture	%	55	52		79		78		54	20		18		19
Anthracene	UG/KG			U	270	J		U.	U		J		Ū	U
Benzaldehyde	UG/KG	89 J		U	110	J	130	J	U	l	J		Ū	U
Benzo(a)Anthracene	UG/KG	U		Ü	150	J		Ü	U		ارز		U	U
Benzo(b)Fluoranthene	UG/KG			U	540	J		U	U	1	J	- 11	Ū	U
Benzo(ghi)Perylene	UG/KG	92 J		U		J		U	U	U	J		ŪΠ	U
Benzo(k)Fluoranthene	UG/KG	U		U	560	J_		U	U		Ĵ		Ü	U
Benzo-a-Pyrene	UG/KG			U	150	J		U	U	l l	J		υ	U
Chrysene	UG/KG	U		Ū	250	J		U	U		J		U	U
Fluoranthene	UG/KG	U		U	340	J		U.	U	L			Ú	U
Indeno (1,2,3-cd) Pyrene	UG/KG	. U		כ	180	J		U	UJ	"   U	J		IJ	U
Naphthalene	UG/KG	U		٦		Ų	3	J	Ü	L	١		J	U
Phenanthrene	UG/KG	U		U	120	J		Ú	U	L	j		J	U
Pyrene	UG/KG	U		U	220	J	120	J	Ü	l l	J		U	U

Ew		-	-	Scan
	THE	ио	1.5	SCHII

	Units	GS01SD		GS02SD	l	GS03SD	GS04SD	L	GS05SD		GS06SD		GS07SD	_ l	GS08SD
2-Methylnaphthalene	UG/KG	0.6	7	0.9	J	33	2	J	0.8	J	0.4	Ĵ	0.4	J	U
Acenaphthene	UG/KG	0.5	J	1	J	100	5		0.9	J		Ū	1	J	Ü
Acenaphthylene	UG/KG	3		2		44	9			Ü		Ü		J	Ü
Anthracene	UG/KG	<b>'</b> 3		7		140	34	L	7			Ū		١	U
Benzo(a)Anthracene	UG/KG	1	J	7		110	19		12			U		J	U
Benzo-a-Pyrene	UG/KG	2		4	Γ.	120	17		16			U		١	υ
Chrysene	UG/KG	1	J	10		200	36	$\Gamma$	26			Ū		٦	U
Dibenzo(a,h)Anthracene	UG/KG		υ	1	J	40	6		. 4		. 1			J	U
Fluoranthene	UG/KG	2		21		190	50		34			Ū	l l	٦٦	U
Fluorene	UG/KG	0.5	J	2		82	6	1	1	J	0.3	J	1	J	U
Naphthalene	UG/KG	0.6	7	. 0.6	J	51	3	J	0.9	J	0.5	J	0.4	1	0.4 J
Phenanthrene	UG/KG	. 2	-	12		72	11	L		U	4		] [	) T	Ū
Pyrene	UG/KG	2		15		140	33		27			U	· ] (	J	Ū

# Miscellaneous Extractables

		GS01SD		GS02SD	GS03SD		GS04SD		GS05SD		GS06SD	GS07SD	GS08SD
.GAMMASITOSTEROL	UG/KG		NR	1400 NJ		NR		NR		NR	NR	NR.	NR
17 UNKNOWN COMPOUNDS	UG/KG		NR	NR		NR		NR		NR	NA	NR	NR
1-DOCOSENE	UG/KG		NR	NR		NR		NR		NR	NR	NR	NR
1-HEXADECENE	UG/KG	540	NJ	NR	1000	NJ		NR		NR	NR	NR	NR
23 UNKNOWN COMPOUNDS	UG/KG	23000	J	NR		NR		NR		NR	NR	NR.	NR
24 UNKNOWN COMPOUNDS	UG/KG		NR	NR		NR	54000	J		NR	NR	NR	NR
25 UNKNOWN COMPOUNDS	UG/KG		NR	NR		NR		NR	140000	J	NR	NR	NR
25 UNKNOWNS	UG/KG		NR	NR	54000	J		NR		ŇR	NR	NR	NR
28 UNIDENTIFIED COMPOUNDS	UG/KG		NR	92000 J		NR		NR		NR	NR	NR	NR
3 UNKNOWN COMPOUNDS	UG/KG		NR	NR		NR		NR		NR	. NR	NR	170 J
BENZENE, 1-METHYL-2-ISOPROPYL	UG/KG !		NR	NR		NR		NR		NJ	NR	NR	NR
BENZO [E] PYRENE	UG/KG		NR	NR	590			NR		NR	NR	NR	NR
CHOLESTANOL	UG/KG		NA	NR		NR	1600	_		NR	NR	NR	NR
ERGOST-5-EN-3-OL, (3.BETA.)-	UG/KG		NR	NR		NR	1700			NR	NR	NR	NR
ERGOST-5-N-3-OL, (3.BETA.)-	UG/KG	1500	NJ	NR.		NR		NR		NR	NR	NR	NR
HEXADECANOIC ACID	UG/KG		NJ	NR		NR		NR		NR	NR	NR	NR
PHENANTHRENONE DERIVATIVE	UG/KG		NR	NR		NR		NR		NJ	NR	NR	NR
STIGMAST-4-EN-3-ONE	UG/KG	730		NR	1400		2200		800		NR	NR	NR
STIGMASTEROL	UG/KG	1100	NJ	NR		NR		NR		NR	NR	NR	NR
UNKNOWN ALCOHOL	UG/KG		NR	NR		NR	9200	J		NR	NR	NR	NR
UNKNOWN CARBOXYLIC ACID	UG/KG		NR	NR		NR		NR	500	J	NR.	NR	NR
UNKNOWN COMPOUND	UG/KG		NR	NR		NR		NR		NR	61 J	69 J	NR
UNKNOWN KETONE	UG/KG		NR	NR		NR	1800	J	•	NR	NR	NR	NR

# Volatiles Scan

	Units :	GS01SD		GS02SD		GS03SD		GS04SD		GS05SD		GS06SD		GS07SD		GS08SD	
% Moisture	%	55		52		79		78		54		20		18		19	
Acetone	UG/KG	94	J	130	J	350	J	520	J	51	J	22 J		49 J		29	J
Carbon Disulfide	UG/KG	5	J		Ū	51	J		U		U	U		U	1		U
Methyl Ethyl Ketone	UG/KG		UJ		W		U	88	J		U	U	IJ	U			UJ

# Miscellaneous Volatiles

		GS01SD	GS02SD	GS03SD	GS04SD		GS05SD	GS06SD	GS07SD	GS08SD
ACETALDEHYDE	UG/KG	NR	NR	NR	71 1	VJ	NR	NR NR	NR	NR
BUTANAL, 3-METHYL-	UG/KG	NR	NR	. NR	1	VR.	NA	NA	NR	NA NA
PROPANAL, 2-MEHTYL-	UG/KG	. NR	NR	NR	1	<b>VR</b>	NR	NR	NR	NR
UNKNOWN COMPOUND	UG/KG	NR	NR	NR	180 J		33 J	NR	NR	8 J

# Page 3 of 3

#### **Data Qualifiers**

U-Analyte not detected at or above reporting limit.

J-Identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Table 5. Groundwater Analytical Results, positive values only. Page 1 of 1

eta		

	Units	GS01GW		GS02GW	[	GS03GW		GS04GW		GS05GW		G\$06GW	
Aluminum	UG/L		υJ	450	J	850	J	1700	J	370	J	74	Γ
Barlum	UG/L	20		120	Ι.	29		26		32			Ū
Beryllium	UG/L	0.11		0.14		0.11			U	0.9			U
Calcium	UG/L	4600		16000		3600		3500		15000			U
Chromlum	UG/L	1.8	R	0.92		1.8		1.6			Ü		U
Cobalt	UG/L	1.2		1.2	R.		U		U	10			U
Copper	UG/L		Ü		Ú		U		U		υ	0.67	
iron	UG/L	1500		1400		2100	Ī	1900		12000			U
Magnesium	UG/L	1500		5500		940		930		8200		,	U
Manganese	UG/L	23		160		110		30		320			U
Nickei	UG/L	1	Ų	3.5			U	3.3		20			U
Potassium	UG/L	520		1500		870		1800		1900		100	
Sodium	UG/L	40000		120000		29000		5000		86000		400	
Vanadium	UG/L		U	0.82	R	. 2		1.9			υ		U
Zinc	UG/L		U	13		18			U.	42			U

#### Pesticides & Aroclors Scan

	GS01GW	GS02GW	GS03GW	GS04GW	GS05GW	GS06GW
No Detected Compounds	1			1.		

#### Extractables Scan

	GS01GW	GS02GW	GS03GW	GS04GW	GS05GW	GS06GW
No Detected Compounds			<del></del>	· · · · · · · · · · · · · · · · · · ·	· · · · · · · · · · · · · · · · · · ·	

#### Miscellaneous Extractables

·		GS01GW		GS02GW		GS03GW		GS04GW	GS05GW		GS06GW	$\Box$
2 UNKNOWN COMPOUNDS	UG/L		NR		NR		NR	NR	4.7	J	NF	A]
D-LIMONENE	UG/L	2.2	NJ		NR		NA	NR		NR	NF	A)
LIMONENE	UG/L		NR	2.2	NJ	2.1	NJ	NR		NR	NF	쥐

### Volatiles Scan

	Units	GS01GW	GS02GW		GS03GW		GS04GW		GS05GW		GS06GW	
Methyl Ethyl Ketone	UG/L	UJ		UJ		UJ	7.2	J		UJ		NR

#### Miscellaneous Volatiles

	Units	GS01GW	GS02GW	GS03GW	GS04GW	GS05GW	GS06GW
INDANE	UG/L_	NR	NR	1.3 NJ	NR.	NR	NR

### **Data Qualifiers**

U-Analyte not detected at or above reporting limit.

J-identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

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L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Table 6. Surface Soll Analytical Results. Page 1 of 10

	Units	GS01SS		GS02SS		GS03SS		GS04SS		GS05SS		GS06SS		GS07SS	GS08SS
% Moisture	%	15		23		25		23		15		26		17	5
Aluminum	MG/KG	4700	J	7800	7	8900	J	4300	J	3100	J	6700	J	2800 J	1800 J
Antimony	MG/KG	0.56	IJ	0.64	R	0.64	U	0.63	U	0.57	U	0.65	Ü	0.58 U	0.5 U
Arsenic	MG/KG	6.3	,	3.3	J	11		2		13		2		3.6	3.4
Barium	MG/KG	72		120		67		60		37		85		34	21
Beryllium	MG/KG	0.33		0.94		0.94		0.37		0.3	Γ.	0.28		0.2	0.19
Cadmium	MG/KG	0.05	٦	0.05	J	0.05	Ü	0.05	5	0.05	U	0.05	J	0.05 U	0.09
Calcium	MG/KG	1100		690		790		690		500		1200		810	800
Chromium	MG/KG	8.4		6.1		20		6		7.5		7.7		4.5	4
Cobalt	MG/KG	5.6		17		8.1	_	1.7		6.6		1.4		2	1.5
Copper	MG/KG	6		. 2	J	6.5	Ü	3.9	U	3	U	4.9	U	3.5 U	2.9 U
Cyanide	MG/KG		NA		NA		NA		NΑ		NA		NA	NA	NA
Iron	MG/KG	10000		8400		27000	J	4500	J	15000	J	5600	J	5400 J	4700 J
Lead	MG/KG	11	J	17		120		18		19		17		11	12
Magnesium	MG/KG	520	J	470	J .	680		280		180		550		250	180
Manganese	MG/KG	630		2800		550		.140		610		170		180	230
Nickel	MG/KG	4.8	U	8.3	Ū	5.1	Ū	2.6	U	2	U	2.8	U	1.9 U	1.7 U
Potassium	MG/KG	380	J	. 240	J	350		200		120		430		170	150
Selenium	MG/KG	1.3		0.93	R	1.9	J	0.63	Ū	0.87	U	0.65	U	0.79 U	0.68 U
Silver	MG/KG	0.39	R	0.71	R	0.63		0.26	R	0.43	_	0.29		0.17 U	0.15 U
Sodium	MG/KG	270		300		82	J	120	J	58	J	110		63	39 U
Thailium	MG/KG	0.73	U	0.81	Ü	0.83	Ū	0.81	Ū	0.73	U	0.84	U	0.75 U	0.65 U
Total Mercury	MG/KG	0.06	U	0.07	U	0.07	υ	0.07	U	0.06	U	0.07	U	0.06 U	0.05 U
Vanadium	MG/KG	17		15		46		11		18		13		8.1	7.1
Zinc	MG/KG	29		47		34		27		23		35		24	26

		GS01SS	GS028S	GS03SS	GS04SS	GS05SS	GS06SS	GS07SS	GS08SS
% Moisture	- 1%	14	23	24	24	15	26	15	4
4,4'-DDD (p,p'-DDD)	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	3.9 U	4.5 U	3.9 U	3.4 U
4,4'-DDE (p,p'-DDE)	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	9.9 U	4.5 U	6.4 U	13 U
4,4'-DDT (p,p'-DDT)	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	4 U	4.5 U	12 J	41 J
Aldrin	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
alpha-BHC	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
alpha-Chlordane /2	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
beta-BHC	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
delta-BHC	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
Dieldrin	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	3.9 U	4.5 U	3.9 U	3.4 U
Endosulfan I (alpha)	UG/KG	2 U	2.2 ∪	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
Endosulfan II (beta)	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	3.9 U	4.5 U	3.9 U	3.4 U
Endosulfan Sulfate	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	3.9 U	4.5 U	3.9 U	3.4 U
Endrin	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	12 U	4.5 U	7.1 U	10 U
Endrin Aldehyde	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	18 J	4.5 U	3.9 U	17 J
Endrin Ketone	UG/KG	3.8 U	4.3 U	4.3 U	4.3 U	17 U	4.5 U	20 U	24 U
gamma-BHC (Lindane)	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
gamma-Chlordane /2	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 0	1.8 U
Heptachlor	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
Heptachlor Epoxide	UG/KG	2 U	2.2 U	2.2 U	2.2 U	2 U	2.3 U	2 U	1.8 U
Methoxychlor	UG/KG	20 U	22 U	22 U	22 U	48 U	23 U	47 U	74 NJ
PCB-1016 (Aroclor 1016)	UG/KG	38 U	43 U	43 U	43 U	39 U	45 U	39 U	34 U
PCB-1221 (Aroclor 1221)	UG/KG	78 U	87 U	88 U	88 U	79 U	91 U	79 U	70 U
PCB-1232 (Aroclor 1232)	UG/KG	38 U	43 U	43 U	43 U	39 U	45 U	39 U	34 U
PCB-1242 (Aroclor 1242)	UG/KG	38 U	43 U	43 U	43 U	39 U	45 U	39 U	34 U
PCB-1248 (Aroclor 1248)	UG/KG	38 U	43 U	43 U	43 U	39 U	45 U	39 U	34 U
PCB-1254 (Aroclor 1254)	UG/KG	38 U	43 U	43 U	43 U	39 U	45 U	39 U	34 U
PCB-1260 (Arocior 1260)	UG/KG	38 U	43 U	43 U	43 U	39 U	45 U	39 U	34 U
Toxaphene	UG/KG	200 U	220 U	220 U	220 U	200 U	230 U	200 U	180 U

		Scan	

	Units	GS01SS	GS02SS	GS035S	GS04SS	GS05SS	GS06SS	GS07SS	GS08SS
% Moisture	%	14	23	24	24	15	26	15	4
(3-and/or 4-)Methylphenol	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
1,1-Biphenyl	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	45 J	1400 U
1,2,4,5-Tetrachlorobenzene	UG/KG	NA	NA NA	NA	NA	NA	NA	NA	NA
2,4,5-Trichlorophenoi	UG/KG	970 U	1100 U	1100 U	1100 U	980 U	1100 U	980 U	3500 U
2,4,6-Trichlorophenol	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2,4-Dichlorophenol	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2,4-Dimethylphenol	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2,4-Dinitrophenol	UG/KG	970 UR	1100 UR	1100 UJ	1100 U	980 U	1100 UJ	980 UJ	3500 UJ
2,4-Dinitrotoluene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2,6-Dinitrotoluene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2-Chloronaphthalene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2-Chiorophenoi	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2-Methyl-4,6-Dinitrophenol	UG/KG	970 U	1100 U	1100 UJ	1100 U	980 U	1100 U	980 UJ	3500 UJ
2-Methylnaphthalene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	120 J	1400 U
2-Methylphenol	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
2-Nitroanline	UG/KG	970 U	1100 U	1100 U	1100 U	980 U	1100 U	980 U	3500 U
2-Nitrophenol	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
3,3'-Dichlorobenzidine	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
3-Nitroaniline	UG/KG	970 U	1100 U	1100 U	1100 U	980 U	1100 U	980 U	3500 U
4-Bromophenyl Phenyl Ether	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
4-Chloro-3-Methylphenol 4-Chloroaniline	UG/KG UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
4-Chlorophenyl Phenyl Ether	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
4-Nitroaniline	UG/KG	970 UJ	1100 UJ	1100 U	1100 U	980 U	1100 U	980 U	3500 U
4-Nitrophenol	UG/KG	970 UJ	1100 UJ	1100 U	1100 U	980 U	1100 U	980 U	3500 U
Acenaphthene	UG/KG	380 U	430 U	430 U	430 Ü	390 U	450 U	170 J	1400 U
Acenaphthylene	UG/KG	380 U	430 U	430 U	430 U	120 J	450 U	320 J	480 J
Acetophenone	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Anthracene	UG/KG	380 U	430 U	430 U	430 U	470	450 U	970	930 J
Atrazine	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Benzaldehyde	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Benzo(a)Anthracene	UG/KG	380 U	430 U	430 U	430 U	770	450 U	2500	1500
Benzo(b)Fluoranthene	UG/KG	380 U	430 U	420 U	430 U	1800 J	450 U	4600 J	11000 J
Benzo(ghi)Perylene	UG/KG	380 U	430 U	430 U	430 U	380 J	450 U	690	3800
Benzo(k)Fluoranthene	UG/KG	380 U	430 U	430 U	430 U	2100 J	450 U	5200 J	12000 J
Benzo-a-Pyrene	UG/KG	380 U	430 U	430 U	430 U	810	450 U	2100	3700
Benzyi Butyi Phthalate	UG/KG	380 U	430 U	430 U	430 UJ	390 UJ	450 U	390 U	1400 U
bis(2-Chloroethoxy)Methane bis(2-Chloroethyl) Ether	UG/KG UG/KG	380 U 380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
bis(2-Chloroisopropyl) Ether	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
bis(2-Ethylhexyl) Phthalate	UG/KG	380 U	430 U	430 U	430 UJ	390 UJ	450 U	390 U	1400 U
Caprolactam	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Carbazole	UG/KG	380 U	430 U	430 U	430 U	200 J	450 U	510	190 J
Chrysene	UG/KG	380 U	430 U	430 U	430 U	1500	450 U	4900	2300
Dibenzo(a,h)Anthracene	UG/KG	380 UJ	430 UJ	430 U	430 U	190 J	450 U	520	1700
Dibenzofuran	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	240 J	1400 U
Diethyl Phthalate	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Dimethyl Phthalate	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
DI-n-Butylphthalate	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Di-n-Octylphthalate	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Fluoranthene	UG/KG	380 U	430 U	430 U	430 U	2600	450 U	8200	2600
Fluorene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	180 J	1400 U
Hexachlorobenzene (HCB)	UG/KG	380 UJ	430 UJ	430 U	430 U	390 U	450 U	390 U	1400 U
Hexachlorobutadiene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Hexachlorocyclopentadiene (HCCP)	UG/KG	380 W	430 UJ	430 U	430 U	390 U	450 UJ	390 U	1400 U
Hexachloroethane	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Indeno (1,2,3-cd) Pyrene	UG/KG	380 UJ	430 UJ	430 U	430 U	610	450 U	1100	5000
Isophorone	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Naphthalene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	87 J	1400 U

Extractables Scan

Volatiles Scan

Toluene
Total Xylenes
trans-1,2-Dichloroethene

Vinyl Chloride

trans-1,3-Dichloropropene
Trichloroethene (Trichloroethylene)
Trichlorofluoromethane (Freon 11)

•	Units	GS01SS	GS0255	GS03SS	GS04SS	GS05SS	G\$06\$\$	GS0/SS	G50855
Nitrobenzene	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
n-Nitrosodi-n-Propylamine	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
n-Nitrosodiphenylamine/Diphenylamine	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Pentachiorophenol	UG/KG	970 UJ	1100 UJ	1100 U	1100 U	980 U	1100 U	980 U	3500 U
Phenanthrene	UG/KG	380 U	430 U	430 U	430 U	320 J	450 U	4200	610 J
Phenol	UG/KG	380 U	430 U	430 U	430 U	390 U	450 U	390 U	1400 U
Pyrene	UG/KG	380 U	430 U	430 U	430 U	1500	450 U	6200	2500

Volatiles Scali	!	GS01SS	Τ	GS02SS	GS03SS		GS04SS		GS05SS		GS06SS	Γ.	GS07SS		GS08SS	7.
% Moisture	%	14		23	24		24		15	_	26		15		· · · · · · · · · · · · · · · · · · ·	NR
1,1,1-Trichloroethane	UQ/KG	11	U	12 U	13	U	12		10 (		13	U	10			NF
1,1,2,2-Tetrachloroethane	UG/KG		Ú	12 U	13		12		10 (			U	10			NF
1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	UG/KG	11	U	12 U	13	U	12		10 (	Ū	13	U	10			NF
1,1,2-Trichloroethane	UG/KG		U	12 U	13		12		10 1	Ú	13	u	10			NF
1,1-Dichloroethane	UG/KG	11	U	12 U	13	U	12	Ü	10 (	U	13	U	10	υT		NF
1,1-Dichloroethene (1,1-Dichloroethylene)	UG/KG	11	U	12 U	13	U	12	U	10	U	13	U	10 1	U		NF
1,2,3-Trichiorobenzene	UG/KG		NA	NA.		NA	Ī .	NA		NA		NA	1	NA		NF
1,2,4-Trichlorobenzene	UG/KG	11	U.	12 U	13	Ū	12	U	10 1	Ū.	13	U	10 (	U		NF
1,2-Dibromo-3-Chloropropane (DBCP)	UG/KG	11	U	12 U	13	IJ		Ū	10	IJ	13	U	10	บป		NF
1,2-Dibromoethane (EDB)	UG/KG	11	Ū	12 U	13	U	12	U	10 (	Ū	13	U	10 1	Ū		NF
1,2-Dichlorobenzene	UG/KG	11	U	12 U	13	U	12		10 (	Ū	13	Ü	10 (	U		NF
1,2-Dichloroethane	UG/KG	11	υ	12 U	13	U	12	C	10 (	J	13	U	10 1	Ū		NF
1,2-Dichloropropane	UG/KG	11	U	12 U	-13	U	12	Ū	10 1	U	13	Ü	10 1	Ū		NF
1,3-Dichlorobenzene	UG/KG	11	U	12 U	13	U	12	v	10 (	Ū	13	U	10	Ū		NF
1,4-Dichlorobenzene	UG/KG	11	U	12 U	13		12	Ü	10 (	Ū.	13	υ	10 (			NF
Acetone	UG/KG	100	J	. 180 J	25	J	28	J	91	J	82	J	87	J		NF
Benzene	UG/KG	11	U	12 U	13	Ü	12		10 1	U	13	Ū	10	Ü		NF
Bromochloromethane	UG/KG	T	NA	\ NA		NA		NA	1	AV		NA		NA		NF
Bromodichloromethane	UG/KG	11	Tu-	12 U	13	C	12	U	10 (	J	13	Ū	10 (	υ		NF
Bromoform	UG/KG	11	U	12 U	13	U	12	حا	10 (	J	13	U	10 (	U		NF
Bromomethane	UG/KG	11	UJ	12 UJ	13	UJ		UJ	10 1	IJ	13	Ū	10 (	UJ		NF.
Carbon Disulfide	UG/KG	11	U	12 U	13	υ	12	U	10 0	J	13	U	10 1	U		NF
Carbon Tetrachioride	UG/KG	11	Ü	12 U	.13	υ	12	c	10 (	Ū	13	Ū	10 (	U		NF
Chlorobenzene	UG/KG	11	U	12 U	13	U	12		10 (	J	13	Ū	10 (			NF
Chloroethane	UG/KG	11	U	12 U	13	Ü	12	U	10 1	J	13	U	10 (	U		NF
Chloroform	UG/KG	11	Ū	12 U	13	Ü	12		10 (	J	13	Ú	10 (	U		NF
Chloromethane	UG/KG	11	Ū	12 U	13	U	12	U	10 (	J	13	Ũ	10 1	Ü		NF
cis-1,2-Dichloroethene	UG/KG	11	U	12 U	13	U	12	U	10 (	J	13	Ū	10 (	Ū		NF
cis-1,3-Dichloropropene	UG/KG	11	U	12 U	13	U	12	U	10 1	J_	13	Ū	10 (	υT		NF
Cyclohexane	UG/KG	11	U	12 U	13	U	12	U	10 L	J	13	U	10 1	U T		NF
Dibromochloromethane	UG/KG	11	U	12 U	13	Ų	12	U	10 1	J	13	Ū	10 (	υT		NF
Dichlorodifluoromethane	UG/KG	11	U	12 U	13	U	12	U	10 (	J	13	U	10 (	Ū İ		NA
Ethyl Benzene	UG/KG	11	U	12 U	13	Ū	12	Ū	10 1	J	13	U	10 (	ΰŤ		NR
Isopropylbenzene	ÚG/KG	11	U	12 U	13	Ū	12	Ū	10 1	j	13	U	10 1	U		NR
Methyl Acetate	UG/KG	11	Ū	12 U	. 13	Ū	12	Ū	10 0	]	13	Ū	10 (	<del>Ú</del>		NR
Methyl Butyl Ketone	UG/KG		Ū	12 U	13		12		10 (		13					NR
Methyl Ethyl Ketone	ÜGÆG	11		21	13		12		10 1		13					NR
Methyl Isobutyl Ketone	UG/KG	11		12 U	13		12		10 L			ŪĴ			<del></del>	NR
Methyl T-Butyl Ether (MTBE)	UG/KG	11	Ū	12 U	13		12		10 L		13		10 0			NE
Methylcyclohexane	UG/KG	11		12 U	13		12		10 1		13		10 0			NR
Methylene Chloride	UG/KG	11		12 U	13		12		10 1		13		10 1			NR
Styrene	UG/KG	11		12 U	13		12		10 0		13		1010			NF
Tetrachloroethene (Tetrachloroethylene)	UG/KG	11		12 0	13		12		10 1		13		10 0			NB
Toluene	UG/KG	11		12 U	13		12		10 0	_	13		10 1			NA
Total Xylenes	UG/KG	11		12 U	13		12		10 0		13		10 0			NR
trans-1.2-Dichloroethene	LIGKG	11		12 U	13		12		10 1			<del>~</del>	101			No

12 U 12 U

12 U

12 U

12 U

13 U

13 U

13 U

13 U

13 U

12 U 12 U 12 U 12 U

12 U

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

11 0

11 U

11 U

11 U

11 U

13 U 13 U

13 U 13 U 13 U

10 U

10 U

10 U

10 U

10 U

10 U

10 U

10 U 10 U

Table 6. Surface Soil Analytical Results. Page 4 of 10

		GS09SS		GS10SS		GS11SS		GS12SS		GS13SS	GS14SS	GS15SS		GS16SS	
% Moisture	%	19		11		7		18		11	14	15		21	
Aluminum	MG/KG	2500	J	4700	J	4400	J	5500	J	4400 J	4800 J	8000	J	7700	J
Antimony	MG/KG	1.2	R	1	R	0.51	บม	0.58	U	0.6 R	0.69 J	0.56	U.	0.7	UJ
Arsenic	MG/KG	0.95	R	6.8		3.7		4.4		6.3	9	6.9		5.5	J
Barium	MG/KG	25		120		54		83		58	65	72		50	
Beryllium	MG/KG	0.15		0.46		0.44		0.46		0.54	0.44	0.71		0.43	
Cadmium	MG/KG	0.05	U	0.34		0.19		0.05	U	0.16 R	0.11 F	0.1		0.05	U
Calcium	MG/KG	340		2600		660		760		1200	1800	2000		1200	
Chromlum	MG/KG	4.1		23		12		8.9		7.8	8.6	11		10	
Cobalt	MG/KG	0.95		4.8		5.4		7.3		4.2	7.3	1.1		. 4	
Copper	MG/KG	0.87	UJ	. 19	J	5	UJ	3.4	Ü	7.2 J	5.7 U	J 4.8	UJ	5.3	Ū
Cyanide	MG/KG		NA		NA		NA		NA	NA	N	A	NA		NA
Iron :	MG/KG	2800		14000		11000		11000	J	12000	11000	12000		12000	
Lead	MG/KG	8.8		80		290		25		33	24	26		15	[J]
Magnesium	MG/KG	160	J	780	J	260	J	380		280 J	340 J	430	J	510	J
Manganese	MG/KG	28		500		880		990		940	870	1800		240	
Nickel	MG/KG	1.5	U	8.5	U	4.9	U	3.2	U	5.2 U	4.1 U	6.7	U	3.6	Ū
Potassium	MG/KG	. 120	J	560	J	330	J	280		390 J	200 J	580	J	490	J
Selenium	MG/KG	0.6	U	0.98	R	0.77	R	1.1	Ū	1.2	0.63 A	1.1	A	1.3	
Silver	MG/KG	0.28	R	0.49		0.38		0.43		0.33 R	0.37	0.5		0.39	R
Sodium	MG/KG	300	$\neg$	320		200		91		230	300	310		270	$\Box$
Thallium	MG/KG	0.77	U	0.7	U	0.66	U	0.75	Ū	0.69 U	0.72 U	0.73	u	0.78	U
Total Mercury	MG/KG	0.06	U	0.06	Ū	0.05	U	0.06	U	0.07 U	0.06 U	0.06	U	0.06	U
Vanadium	MG/KG	6.6		18		16		22		17	21	23		22	
Zinc	MG/KG	19	J	160		840		28		72	44	48		33	

		GS09SS	GS10SS	GS11SS	GS12SS	G\$13SS	GS14SS	GS15SS	GS16SS
% Moisture	.%	26	12	7	17	13	14	16	21
4,4'-DDD (p,p'-DDD)	UG/KG	4.5 U	3.8 U	3.5 U	4 U	3.8 U	3.8 U	3.9 U	4.2 U
4,4'-DDE (p,p'-DDE)	UG/KG	4.5 U	3.8 ∪	3.5 U	4 U	30 U	3.8 U	3.9 U	2.5 J
4,4'-DDT (p,p'-DDT)	UG/KG	4.5 U	43 J	30 J	4 U	38 J	10 NJ	30 NJ	4.2 U
Aldrin	UG/KG	2.3 U	1.9 ∪	1.8 U	2 U	2 U	2 U	2 U	2.2 U
alpha-BHC	UG/KG	2.3 U	1.9 U	1.8 U	2 U	2 U	2 U	2 U	2.2 U
alpha-Chiordane /2	UG/KG	2.3 U	1.9 U	1.8 U	2 U	2 U	2 U	2 U	2.2 U_
beta-BHC	UG/KG	2.3 U	12 U	5.9 U	2 U	2 U	2 U	2 U	2.2 U
deita-BHC	UG/KG	2.3 U	1.9 U	1.8 U	2 U	2 U	2 U	2 U	2.2 U
Dieldrin	UG/KG	4.5 U	3.8 U	3.5 U	4 U	3.8 U	3.8 U	3.9 U	4.2 U
Endosullan I (alpha)	UG/KG	2.3 U	1.9 U	1.8 U	2 U	2 U	2 U	2 U	2.2 U
Endosulfan II (beta)	UG/KG	4.5 U	3.8 U	3.5 U	4 U	3.8 U	3.8 U	3.9 U	4.2 U
Endosulfan Sulfate	UG/KG	4.5 U	3.8 U	3.5 U	4 U	3.8 U	3.8 U	3.9 U	4.2 U
Endrin	UG/KG	4.5 U	15 U	3.5 U	4 U	42 J	3.8 U	11 U	4.2 U
Endrin Aldehyde	UG/KG	4.5 U	3.8 U	3.5 U	4 U	49 J	3.8 U	27 NJ	4.2 U
Endrin Ketone	UG/KG	4.5 U	3.8 U	3.5 U	4 U	150	3.8 U	3.9 U	14 J
gamma-BHC (Lindane)	UG/KG	2.3 ∪	1.9 U	1.8 U	2 U	2 0	2 U	2 U	2.2 U
gamma-Chlordane /2	UG/KG	2.3 U	1.9 U	1.8 U	2 U	18 U	2 U	2 U	2.2 U
Heptachlor	UG/KG	2.3 U	1.9 U	1.8 U	2 0	2 U	2 U	2 U	2.2 U
Heptachlor Epoxide	UG/KG	2.3 U	1.9 U	1.8 U	2 U	2 U	2 U	2 U	2.2 U
Methoxychlor	UG/KG	23 U	81 NJ	54 U	20 U	20 U	34 NJ	61 U	22 U
PCB-1016 (Aroclor 1016)	UG/KG	45 U	38 U	35 U	40 U	38 U	38 U	39 U	42 U
PCB-1221 (Aroclor 1221)	UG/KG	91 U	76 U	72 U	81 U	77 U	78 U	80 U	85 U
PCB-1232 (Aroclor 1232)	UG/KG_	45 U	38 U	35 U	40 U	38 U	38 U	39 U	42 U
PCB-1242 (Arocior 1242)	UG/KG	45 U	38 U	35 U	40 U	38 U	38 U	39 U	42 U
PCB-1248 (Arocior 1248)	UG/KG	45 U	38 U	35 U	40 U	38 U	38 U	39 U	42 U
PCB-1254 (Aroclor 1254)	UG/KG	45 U	38 U	35 U	40 U	38 U	38 U	39 U	42 U
PCB-1260 (Aroclor 1260)	UG/KG	45 U	38 U	35 U	40 U	38 U	38 U	39 U	42 U
Toxaphane	UG/KG	230 U	190 U	180 Ú	200 U	200 U	200 U	200 U	220 U

Ext	ra	cta	ble	26	Sc	an

EXTROMOTO COUNT	Units	GS09SS	GS10SS	GS11SS	GS12SS	GS13SS	GS14SS	GS15SS	GS16SS
% Moisture	%	26	12	7	17	13	14	16	21
(3-and/or 4-)Methylphenol	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
1,1-Biphenyl	UGAKG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
1,2,4,5-Tetrachiorobenzene	UG/KG	NA	NA	NA	NA NA	NA NA	. NA	NA	NA
2,4,5-Trichlorophenol	UG/KG	1100 U	940 U	890 U	1000 U	2900 U	970 U	990 U	1100 U
2,4,6-Trichlorophenol	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
2,4-Dichlorophenol	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
2,4-Dimethylphenoi	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
2,4-Dinitrophenol	UG/KG	1100 UJ	940 UJ	890 UJ	1000 UJ	2900 U	970 UJ	990 UR	1100 U
2,4-Dinitrotoluene	UG/KG	450 U	380 U	350 U	400 U	1100 U	. 380 U	390 U	420 U_
2,6-Dinitrotoluene	UG/KG	· 450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
2-Chloronaphthalene	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
2-Chlorophenol	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
2-Methyl-4,6-Dinitrophenol	UG/KG	1100 U	940 U	890 U	1000 UJ	2900 U	970 U	990 U	1100 U
2-Methylnaphthalene	nd/kg	450 U	69 J	70 J	400 U	1100 U	380 U	390 U	420 U
2-Methylphenol	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
2-Nitroanliine	UG/KG	1100 U	940 U	890 U	1000 U	2900 U	970 U	990 U	1100 U
2-Nitrophenol	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
3,3'-Dichlorobenzidine	UG/KG	450 U	380 U	350 U	400 U	1100 UJ	380 U	390 U	420 UJ
3-Nitroaniline	UG/KG	1100 U	940 U	890 U	1000 U	2900 U	970 U	990 U	1100 U
4-Bromophenyl Phenyl Ether	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
4-Chloro-3-Methylphenoi	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
4-Chlorophand Phand Ethan	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
4-Chlorophenyl Phenyl Ether 4-Nitroaniline	UG/KG UG/KG	1100 U	380 U	350 U 890 U	1000 U	1100 U 2900 U	380 U 970 U	390 U 990 UJ	1100 U
4-Nitrophenol	UG/KG	1100 U	940 U	890 U	1000 U	2900 U	970 U	990 UJ	1100 U
Acenaphthene	UG/KG	450 U	97 J	350 U	400 U	1100 U	380 U	390 U	420 U
Acenaphthylene	UG/KG	450 U	1100	330 J	400 U	1000 J	380 U	250 J	87 J
Acetophenone	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Anthracene	UG/KG	450 U	1700	830	400 U	1600	1100	370 J	420 U
Atrazine	UG/KG	450 U	380 U	350 U	400 Ü	1100 U	380 U	390 U	420 U
Benzaldehyde	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Benzo(a)Anthracene	UG/KG	450 U	2800	1000	400 U	4300	2300	690	260 J
Benzo(b)Fluoranthene	UG/KG	450 U	5200	2500	400 U	7300 J	4200	1500	700 J
Benzo(ghl)Perylene	UG/KG	450 U	1300	460	400 U	2000	820	360 J	92 J
Benzo(k)Fluoranthene	UG/KG	450 U	2300	1700	400 U	5400 J	2400	1100	730 J
Benzo-a-Pyrene	UG/KG	450 U	3100	800	400 U	3100	2300	650	230 J
Benzyl Butyl Phthalate	UG/KG	450 U	380 U	350 U	400 U	1100 UJ	380 U	390 U	420 UJ
bis(2-Chloroethoxy)Methane	UG/KG	450 U	380 U	350 U	· 400 U	1100 U	380 U	390 U	420 U
bis(2-Chloroethyl) Ether	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
bis(2-Chiorolsopropyl) Ether	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
bis(2-Ethylhexyl) Phthalate	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Caprolactam	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Carbazole	UG/KG	450 U	480	210 J	400 U	540 J	280 J	250 J	62 J
Chrysene	UG/KG	450 Ü	4300	1500	400 U	6200	3100	1700	390 J
Dibenzo(a,h)Anthracene	UG/KG	450 U	750	310 J	400 U	1100	580	200 J	95 J
Dibenzofuran	UG/KG UG/KG	450 U 450 U	120 J 380 U	140 J 350 U	400 U	1100 U	380 U	390 U	420 U
Olethyl Phthalate Dimethyl Phthalate	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
	UG/KG	450 U	380 U	350 U	400 U	1100 U		390 U	420 U
Di-n-Butylphthalate Di-n-Octylphthalate	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Fluoranthene	UG/KG UG/KG	450 U	4700	1900	400 U	11000	3100	3100	760
Fluorene	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Hexachlorobenzene (HCB)	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 UJ	420 U
Hexachlorobutadione	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Hexachiorocyclopentadiene (HCCP)	UG/KG	450 U	380 UJ	350 U	400 U	1100 U	380 U	390 UJ	420 U
Hexachioroethane	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Indeno (1,2,3-cd) Pyrene	UG/KG	450 U	2300	940	400 U	3100	1700	740 J	270 J
Isophorone	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Naphthalene	UG/KG	450 U	110 J	120 J	400 U	1100 U	380 U	390 U	420 U
(ABAILTIGIOID	JUNE	40010	11010	12013	70010	110010	30010	330 0	720 0

# Extractables Scan

	Units	GS09SS	GS10SS	GS11SS	GS12SS	GS13SS	GS14SS	GS15SS	GS16SS
Nitrobenzene	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
n-Nitrosodi-n-Propylamine	UG/KG	450 U	380 U	350 U	400 U	1100 Ü	380 U	390 U	420 U
n-Nitrosodiphenylamine/Diphenylamine	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Pentachlorophenol	UG/KG	1100 U	690 J	640 J	1000 U	2900 U	970 U	990 UJ	1100 U
Phenanthrene	UG/KG	450 U	1400	540	400 U	1600	320 J	1100	310 J
Phenol	UG/KG	450 U	380 U	350 U	400 U	1100 U	380 U	390 U	420 U
Pyrene	UG/KG	450 U	5000	1700	400 U	6800	3300	2200	540

### Volatiles Scan

		GS09SS	GS10SS	GS11SS	GS12SS	GS13SS	GS14SS	GS15SS	GS16SS
% Moisture	.%	26	12	7	17	13	14	16	21
1,1,1-Trichioroethane	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 0	10 U	12 U
	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	UG/KG	13 U	11 U	11 (	10 U	15 U	11 U	10 U	12 U
1,1,2-Trichloroethane	UG/KG	13 U	11 U	11 1	J 10 U	15 U	11 U	10 U	12 U
1,1-Dichioroethane	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
1,1-Dichloroethene (1,1-Dichloroethylene)	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
1,2,3-Trichlorobenzene	UG/KG	NA	N/	A	NA NA	A NA	NA	. NA	NA NA
1,2,4-Trichlorobenzene	UG/KG	. 13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
1,2-Dibromo-3-Chioropropane (DBCP)	UG/KG	13 U	11 U	11 (	J 10 U	15 UJ	11 U	10 U	12 U
1,2-Dibromoethane (EDB)	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
1,2-Dichlorobenzene	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
1,2-Dichloroethane	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
1,2-Dichloropropane	UG/KG	13 U	11 U	11 1	J 10 U	15 U	11 U	10 U	12 U
1,3-Dichlorobenzene	UG/KG	13 U	11 U	11 (	J 10 U	15 U	11 U	10 U	12 U
17	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 UJ	84 J	220 J			75 J	100 J	56 J
	UG/KG	13 U	11 U	11 (		15 U	11 U	10 U	12 U
Bromochloromethane	UG/KG	· NA	N/	N	IA NA	NA NA	NA NA		
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11_0	10 U	12 U
	UG/KG	13 UJ	11 U.			15 UJ	11 UJ	10 UJ	12 UJ
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	ug/kg	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 (		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	UG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 L		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 6		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	3 1	12 U
	JG/KG	13 UJ	11 U	11 U			11 U	10 U	12 U
	JG/KG	13 UJ	11 U	21 J		15 J	11 U	12	12 U
	JG/KG	13 UJ	11 U	11 U			11 U	10 U	12 U
	JG/KG	13 UJ	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
	JG/KG	13 U	11 U	11 U		15 U	11 U	10 U	12 U
Vinyi Chloride	JG/KG	13 U	11 U	11 U	10 U	15 U	11 U	10 U	12 U

Table 6. Surface Soll Analytical Results. Page 7 of 10

	<u> </u>	GS17SS	GS18SS	GS19SS	GS20SS	GS20SD	GS21SS	GS22SS	GS23SS	G\$24\$\$
% Moisture	%	21	22	15	20	19	22	14	26	47
Aluminum	MG/KG	9000 J	4900 J	8500 J	6700 J	6100 J	6800 J	4200 J	7500 J	4900 J
Antimony	MG/KG	0.83 U.	1 00	0.89 UJ	0.6 UJ	0.59 UJ	0.96 UJ	0.56 U	0.65 UJ	0.91 U
Arsenic	MG/KG	5.5	5.4	6.5	5.8	4.5	2.7	4	5	6.5
Baríum	MG/KG	120	53	89	410	130	88	42	83	120
Beryllium	MG/KG	0.94	0.35	0.73	1.2	0.76	0.81	0.43	0.5	0.61
Cadmium	MG/KG	0.07 R	0.05 U	0.05 U	0.17	0.05 U	0.05 U	0.05 U	0.1 R	0.16
Calcium	MG/KG	2300	890	2000	1700	1400	190	280	960	1600
Chromlum	MG/KG	8	9.9	8.3	8.2	7.2	5.6	9.4	12	12
Cobalt	MG/KG	12	7.6	13	17	6.8	14	5.5	6.8	6.8
Copper	MG/KG	5.5 U	3.1 U	3.8 U	0.95 U	4.3 U	1.5 U	2.5 U	12 J	8.4 U
Cyanide	MG/KG	N/	A NA	NA.	NA.	NA	NA	NA	NA	NA NA
Iron	MG/KG	12000	10000	12000	12000	8800	7300	9700 J	13000	17000 J
Lead	MG/KG	20 J	15 J	19 J	24 J	22 J	13 J	9.8	24	28
Magnesium	MG/KG	490 J	340 J	530 J	380 J	410 J	270 J	220	630 J	450
Manganese	MG/KG	2500	590	1900	5500	1800	2200	470	920	890
Nickel	MG/KG	9.2 U	3.9 U	6.7 U	8 U	6.1 U	6.7 U	2.2 U	4.8 U	5.3 U
Potassium	MG/KG	620 J	180 J	550 J	240 J	250 J	250 J	170	480 J	370
Selenium	MG/KG	1.5 J	1.1	1.2	1.3 J	1.1 R	0.8	1.1 U.	1.1 R	1.5 U
Silver	MG/KG	0.51 R	0.29	0.53 R	0.9 R	0.42	0.52	0.2	0.46 R	0.47
Sodium	MG/KG	280	310	290	. 290	290	300	44 U	340	86 J
Thaillum	MG/KG	0.79 U	0.8 U	0.73 U	1.5	0.77 U	0.79 U	0.72 U	0.83 U	1.2 U
Total Mercury	MG/KG	0.08 U	0.06 U	0.07 U	0.08 U	0.07 U	0.06 U	0.06 U	0.08 U	0.09 U
Vanadium	MG/KG	23	20	22	22	17	13	21	23	26
Zinc	MG/KG	65	34	52	88	63	24	19	49	89

			GS18SS	GS19SS	GS20SS	GS20SD	GS21SS	GS22SS	GS23SS	GS24\$S
% Moisture	%	GS17\$S	23	17	21	20	22	13	26	45
4,4'-DDD (p,p'-DDD)	UG/KG	4.2 U	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	4.5 U	6 U
4,4'-DDE (p,p'-DDE)	UG/KG	4.2 U	4.3 U	4 U	4.1 NJ	3.3 NJ	4.2 U	.3.8 U	4.5 U	6 U
4,4'-DDT (p,p'-DDT)	UG/KG	4.2 U	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	4.5 U	6 U
Aldrin	UG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	2.3 U	3.1 U
alpha-BHC	UG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	1.7 NJ	3.1 U
alpha-Chlordane /2	UG/KG	2.2 U	2.2 U	2 U	-2.2 U	2.1 U	2.2 U	2 U	2.3 U	3.1 U
beta-BHC	UG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	2.3 U	5.3 U
delta-BHC	UG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	2.3 U	3.1 U
Dieidrin	UG/KG	4.2 U	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	4.5 U	6 U
Endosulfan i (aipha)	UG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	2.3 U	3.1 U
Endosulfan II (beta)	UG/KG	4.2 U	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	4.5 U	6 U
Endosultan Sulfate	UG/KG	4.2 U	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	4.5 U	6 U
Endrin	UG/KG	6.4 J	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	4.5 U	6 U
Endrin Aldehyde	UG/KG	4.2 U	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	4.5 U	. 6 U
	UG/KG	20 J	4.3 U	4 U	4.2 U	4.1 U	4.2 U	3.8 U	6.9 J	6 U
gamma-BHC (Lindane)	₩G/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	2.3 U	3.1 U
gamma-Chlordane /2	ΨG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	2.3 ↓	3.1 U
Heptachlor	ΨG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U	2 U	2.3 U	3.1 U
Heptachior Epoxide	UG/KG	2.2 U	2.2 U	2 U	2.2 U	2.1 U	2.2 U_	2 U	2.3 U_	3.1 U
Methoxychlor	UG/KG	22 U	22 U	20 U	22 U	21 U	22 U	20 U	23 U	31 U
PCB-1016 (Arocior 1016)	UG/KG	42 U	43 U	40 U	42 U	41 U	42 U	38 U	45 U	60 U
	UG/KG	85 U	87 U	81 U	85 U	84 U	86 U	77 U	91 U_	120 U
PCB-1232 (Aroclor 1232)	UG/KG	42 U	43 U	40 U	42 U	41 U	42 U	38 U	45 U	60 U
PCB-1242 (Aroclor 1242)	UG/KG	42 U	43 U	40 U	42 U	41 U	42 U	38 U	45 U	60 U
PCB-1248 (Aroclor 1248)	UG/KG	42 U	43 U	40 U	42 U	41 U	.42 U	38 U	45 U	60 U
	UG/KG	42 U	43 U	40 U	42 U	41 U	42 U	38 U	45 U	60 U
PCB-1260 (Aroclor 1260)	UG/KG	42 U	43 U	40 U	42 U	41 U	42 U	38 U	45 U	60 U
Toxaphene	UG/KG	220 U	220 U .	200 U	220 U	210 U	220 U	200 U	230 U	310 U

	Units	GS17SS	GS18SS	GS19SS	GS20SS	GS20SD	GS21SS	GS22SS	GS23SS	GS24SS
% Moisture	*	21-	23	17	21	20	22	13	26	45
(3-and/or 4-)Methylpheno!	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
1,1-Bipheny!	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
1,2,4,5-Tetrachlorobenzene	UG/KG	. NA	NA	NA	NA	NA	NA	NA	NA	NA
2,4,5-Trichlorophenol	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 U	1100 U	1500 U
2,4,6-Trichlorophenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2,4-Dichlorophenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2,4-Dimethylphenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2,4-Dinitrophenol	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 UJ	1100 UR	1500 UJ
2,4-Dinitrotoluene	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2,6-Dinitrotoluene	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2-Chloronaphthalene	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2-Chlorophenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2-Methyl-4,6-Dinitrophenol	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 U	1100 U	1500 UJ
2-Methylnaphthaiene	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2-Methylphenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
2-Nitroaniline	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 U	1100 U	1500 U
2-Nitrophenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
3.3'-Dichlorobenzidine	UG/KG	420 UJ	430 U	400 U	420 UJ	410 UJ	420 UJ	380 U	450 U	600 U
3-Nitroaniline	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 U	1100 U	1500 U
4-Bromophenyl Phenyl Ether	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U.
4-Chloro-3-Methylphenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
4-Chloroaniline	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	000 U
4-Chlorophenyl Phenyl Ether	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
4-Nitroanline	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 U	1100 UJ	1500 U
4-Nitrophenol	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 U	1100 UJ	1500 U
Acenaphthene	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Acenaphthylene	UG/KG	270 J	430 U	400 U	420 U	410 U	420 U	380 U	66 J	150 J
Acetophenone	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Anthracene	UG/KG	380 J	430 U	400 U	420 U	410 U	420 U	380 U	430 U	330 J
Atrazine	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Benzaldehyde	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Benzo(a)Anthracene	UG/KG	1000	430 U	400 U	58 J	44 J	420 U	380 U	430 J	990
Benzo(b)Fluoranthene	UG/KG	2600 J	430 U	400 U	160 J	140 J	420 U	380 U	NA	2700 J
Benzo(ghi)Perylene	UG/KG	290 J	430 U	400 U	420 U	410 U	420 U	380 U	180 J	330 J
Benzo(k)Fluoranthene	UG/KG	2700 J	430 U	400 U	170 J	150 J	420 U	380 U	520	2600 J
Benzo-a-Pyrene	UG/KG	1000	430 U	400 U	46 J	47 J	420 U	380 U	380 J	870
Benzyl Butyl Phthalate	UG/KG	420 UJ	430 UJ	400 UJ	420 UJ	410 UJ	420 UJ	380 U	450 U	600 U
bis(2-Chloroethoxy)Methane	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
bis(2-Chioroethyi) Ether	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
bis(2-Chloroisopropyl) Ether	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
bis(2-Ethylhexyl) Phthalate	ÜG/KG	420 U	430 UJ	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Caprolactam	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Carbazole	UG/KG	130 J	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Chrysene	UG/KG	1600	430 U	400 U	98 J	69 J	420 U	380 U	510	1100
Dibenzo(a,h)Anthracene	UG/KG	330 J	430 U	400 U	420 U	410 U	420 U	380 U	120 J	260 J
Dibenzofuran	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Diethyl Phthalate	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Dimethyl Phthalate	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Di-n-Butyiphthalate	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Di-n-Octylphthalate	UG/KG	420 U	430 U	400 U	420 U	410 U	420 Ü	380 U	450 U	600 U
Fluoranthene	UG/KG	2500	430 U	400 U	73 3	81 J	420 U	380 U	520	1300
Fluorene	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Hexachlorobenzene (HCB)	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 UJ	600 U
Hexachlorobutadiene	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Hexachiorocyclopentadiene (HCCP)	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U		450 UJ	600 U
Hexachioroethane	UG/KG	420 U	430 U	400 U				380 U	450 U	600 U
Indeno (1,2,3-cd) Pyrene	UG/KG	890			. 420 U	410 U	420 U	380 U		
			430 U	400 U	63 J	51 J	420 U	380 U	340 J	550 J
Isophorone	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	600 U
Naphthalene	:UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	. 450 U	600 U

### Extractables Scan

	Units	GS17SS	GS18SS	GS19SS	G\$20\$\$	GS20SD	G\$2155	GS22SS	GS23SS	GS2455
Nitrobenzene	UG/KG	420 U	430 U	400 U	420 L	J 410 U	420 U	380 U	450 U	600 U
n-Nitrosodi-n-Propylamine	UG/KG	420 U	430 U	400 U	420 L	410 U	420 U	380 U	450 U	600 U
n-Nitrosodiphenylamine/Diphenylamine	UG/KG	420 U	430 U	400 U	420 L		420 U	380 U	450 U	600 U
Pentachlorophenoi	UG/KG	1100 U	1100 U	1000 U	1100 U	1000 U	1100 U	950 U	1100 UJ	1500 U
Phenanthrene	UG/KG	250 J	430 U	400 U	420 U	/410 U	420 U	380 U	59 J	600 U
Phenol	UG/KG	420 U	430 U	400 U	420 U	410 U	420 U	380 U	450 U	_ 600 U
Pyrene	UG/KG	1700	430 U	400 U	100 J	63 J	420 U	380 U	450	1600

Moisture  1,1,1-Trichloroethane 1,1,2,2-Tetrachloroethane 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 1,1,2-Trichloroethane 1,1-Dichloroethane 1,1-Dichloroethane 1,1-Dichloroethane 1,2,3-Trichlorobenzene 1,2-Dibromo-3-Chloropropane (DBCP) 1,2-Dibromo-3-Chloropropane (DBCP) 1,2-Dibromoethane (EDB) 1,2-Dichlorobenzene	% UGMG UGMG UGMG UGMG UGMG UGMG	951755 21 13 U 13 U 13 U 13 U 13 U	J J	23 13 U 13 U 13 U	GS19SS 17 11	11	GS20SS 21	GS20SD 20	GS21SS 22	GS22SS 13	GS23SS	GS24SS
,1,1-Trichloroethane ,1,2,2-Tetrachloroethane ,1,2,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) ,1,2-Trichloroethane ,1-Dichloroethane ,1-Dichloroethane ,1-Dichloroethane (1,1-Dichloroethylene) ,2,3-Trichlorobenzene ,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	13 U 13 U 13 U 13 U 13 U	) ) )	13 U 13 U	11	11		1 201	221			
,1,2,2-Tetrachioroethane ,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) ,1,2-Trichloroethane ,1-Dichloroethane ,1-Dichloroethane (1,1-Dichloroethylene) ,2,3-Trichlorobenzene ,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene	UG/KG UG/KG UG/KG UG/KG	13 L 13 L 13 L	) ) )	13 U			40.11	44 11		11 0	26	45 25 U
,1,2-Trichloro-1,2,2-Trifliuoroethane (Freon 113) ,1,2-Trichloroethane ,1-Dichloroethane ,1-Dichloroethane (1,1-Dichloroethylene) ,2,3-Trichlorobenzene ,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene	UG/KG UG/KG UG/KG	13 L 13 L 13 L	J				12 U	11 U			13 U	
,1,2-Trichloroethane ,1-Dichloroethane ,1-Dichloroethene (1,1-Dichloroethylene) ,2,3-Trichlorobenzene ,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene	UG/KG UG/KG UG/KG	13 U	Л	1310			12 U	11 U	12 U	11 U	13 U	25 U
,1-Dichloroethane ,1-Dichloroethene (1,1-Dichloroethylene) ,2,3-Trichlorobenzene ,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene	UG/KG UG/KG	13 L			11		12 U	11 0	12 U	11 U	13 U	25 U
,1-Dichloroethene (1,1-Dichloroethylene) ,2,3-Trichlorobenzene ,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene	UG/KG			13 U	11		12 U	11 U		11 U	13 U	25 U
,2,3-Trichlorobenzene ,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene		401		13 U	11		12 U	11 U	12 U		13 U	25 U
,2,4-Trichlorobenzene ,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene		13 L		13 U	11		12 U	11 U		11 U	13 U	. 25 U
,2-Dibromo-3-Chloropropane (DBCP) ,2-Dibromoethane (EDB) ,2-Dichlorobenzene			IA	NA NA		NΑ	NA 10					
,2-Dibromoethane (EDB) ,2-Dichlorobenzene	UG/KG	13 U		13 U	11		12 U	11 U		11 U	13 U	25 U
,2-Dichlorobenzene	UG/KG	13 6		13 UJ	11		12 U	11 U.	12 UJ 12 U		13 U	25 U
	UG/KG	13 L		13 U	11		12 U	11 U		11 U	13 U	25 U
O Diables athers	UG/KG UG/KG	13 L		13 U	11		12 U	11 U	12 U	11 U	13 U	25 U
,2-Dichloroethane		13 0			.11		12 U	11 0	12 U	11 U		25 U
,2-Dichioropropane ,3-Dichiorobenzane	UG/KG UG/KG	13 0		13 U	11			11 0	12 U	11 0	13 U	25 U
,3-Dichlorobenzene	UG/KG	13 0		13 U	11		12 U 12 U	11 0	12 U	11 0	13 U	25 U
Acetone	UG/KG	260 J		130 J	180		83 J	230 J	270 J	110 J	130 J	25 U
Benzene	UG/KG	13 U		13 U	11		12 U	11 U	12 U	11 U	13 U	25 U
Promochloromethane	UG/KG		IA -	NA NA		NA.	NA					
Promodichloromethane	UG/KG	13 0		13 U	11		12 U	11 0	12 U	11 0	13 U	25 U
Bromoform	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 0	25 U
Promomethane	UG/KG	13 0		13 UJ	11		12 UJ	11 0.			13 0	
Carbon Disuifide	UG/KG	13 0		13 U	11		12 U	11 0	12 U	11 0	13 0	25 0
Carbon Tetrachloride	UG/KG	13 0		13 U	11		12 U	11 0	12 U	11 0	13 U	25 U
Chlorobenzene	UG/KG	13 U		13 0	11		12 U	11 0	12 U	11 0	13 U	25 U
Chloroethane	UG/KG	13 U		13 U	11		12 U	11 0	12 U	1110	13 0	25 U
hioroform	UG/KG	13 U		13 0	• 11		12 U	1110	12 U	11 0	13 U	25 U
Chloromethane	UG/KG	13 U		13 U	11		12 U	11 0	12 U	1110	13 U	25 U
is-1,2-Dichloroethene	UG/KG	13 U		13 U	11		12 U	1110	12 U	11 0	13 U	25 U
ls-1,3-Dichloropropene	UG/KG	13 0		13 0	11		12 U	1110	12 U	11 0	13 0	25 U
cyclohexane	UG/KG	13 U		13 U	11		12 U	1110	12 U	11 0	13 U	25 U
Dibromochloromethane	UG/KG	13 U		13 U	11		12 U	11 0	12 U	110	13 U	25 U
Dichlorodiffuoromethane	UG/KG	13 U		13 U	11		12 U	11 0	12 U	1110	13 U	25 U
thyl Benzene	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 U	25 U
topropylbenzene	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 U	25 U
lethyl Acetate	UG/KG	13 U		13 U	3		12 U	1110	12 U	11 0	13 U	25 U
lethyl Butyl Ketone	UG/KG	13 U		13 UJ	11		12 U	11 0		11 03		25 U
lethyl Ethyl Ketone	UG/KG	28 J		14 J	17		12 U	23 J	31 J	11 03		25 UJ
lethyl Isobutyl Ketone	UG/KG	13 U		13 00	11		12 U	11 03		11 03	13 U	25 UJ
lethyl T-Butyl Ether (MTBE)	UG/KG	13 U		13 0			12 U	11 0	12 U	11 0	13 U	25 U
lethylcyclohexane	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 U	25 U
lethylene Chloride	UG/KG	13 U		13 0	11		12 U	11 0	12 U	11 0	13 U	25 U
tyrene	UG/KG	13 U		13 U	11/0		12 U	11 U	12 U	1110	13 0	25 U
strachloroethene (Tetrachloroethylene)	UG/KG	13 U		13 0	11		12 U	11 0	12 U	11 0	13 U	25 U
oluene	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 U	25 U
otal Xylenes	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 0	25 U
ans-1,2-Dichloroethene	UG/KG	13 U		13 0	11		12 U	11 0	12 0	11 0	13 U	25 U
ans-1,3-Dichloropropene	UG/KG	13 U		13 0	11		12 U	11 0	12 U	11 0	13 0	25 U
richloroethene (Trichloroethylene)	UG/KG	13 U		13 0	11		12 U	11 0	12 U	11 0	13 0	25 0
richiorofluoromethane (Freon 11)	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 U	25 U
Inyl Chloride	UG/KG	13 U		13 U	11		12 U	11 0	12 U	11 0	13 0	25 U

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#### **Data Qualifiers**

U-Analyte not detected at or above reporting limit.

J-identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Table 7. Subsurface Soil Analytical Results. Page 1 of 10

•	i	GS01SB		GS02SB	GS03SB		GS04SB	Γ	GS05SB	GS06SB	GS07SB		GS08SB	
% Moisture	%	18		19	28		27		23	27	26		18	
Aluminum	MG/KG	7300	7	7800 J	8600	J	6600	J_	5300 J	7700 J	9400	J	8100	
Antimony	MG/KG	0.59	C)	0.59 UJ	0.67	C	0.66	U	0.62 U	0.65 U	0.65	U	0.58	
Arsenic	MG/KG	6.7		3.4 J	9.4		1.9		0.91 U	2.7	2.7		3.5	
Barium	MG/KG	62		34	63		57		59	48	50		45	
Beryllium	MG/KG	0.31		0.36	0.64		0.41		0.27	0.24	0.35		0.59	
Cadmium	MG/KG	0.05	Ü	0.05 U	0.06	U	0.06	ح	0.05 U	0.05 U	0.05	U	0.05	Ū
Calcium	MG/KG	850		170	1400		670		580	990	670		450	
Chromium	MG/KG	12		9.5	14		6.6		4.9	7.5	8	T	9.2	
Cobalt	MG/KG	3.3		2.6	4.3		1.4		0.78 R	1.2	1.6		9.4	
Copper	MG/KG	6.4		3.8 UJ	6	Ū.		ح	2.1 U	3.4 U	5.7 (	Ų T	3.9	U
Cyanide	MG/KG		NA	NA		NA		Š	NA NA	NA	1	NA		NA
Iron	MG/KG	13000		13000	22000	_	7500	J	2300 J	5800 J	8200	J	12000	J
Lead	MG/KG	8.2	J	6.6	15		9.2		. 6.7	11	11	T	5.7	
Magnesium	MG/KG	580	J	480 J	770		370		220	580	550		370	
Manganese	MG/KG	180		200	z 200		60		16	18	19	1	1000	
Nickel	MG/KG	5.2	U	4.1 U	5.1	C		Ü	1.7 U	2.7 U	3.7 l	U	5.3	U
Potassium	MG/KG	_320	J	320 J	360		190		170	240	260		350	
Selenium	MG/KG	0.7		1.4	1.6		0.66	C	0.62 U	0.96 U	0.65	U_T	0.58	U
Silver	MG/KG	0.38		0.38 R	0.49	R	0.19	U	0.18 U	0.28 R	0.19 (	J	0.32	
Sodium	MG/KG	290		290	98	5	300		190 J	250	220		46	
Thallium	MG/KG	0.76	Ū	0.77 U	0.87	Ú	0.85		0.81 U	0.84 U	0.84 (	<u> </u>	0.75	Ū
Total Mercury	MG/KG	0.06	Ü	0.06 U	0.07	Ü	0.07	U	0.07 U	0.07 U	0.07 l	U	0.06	U
Vanadium	MG/KG	22		20	33		14		6.7	17	18	$\Box \Box$	22	
Zinc	MG/KG	18		20	23		15		7.3	12	15	$\perp I$	25	

	1.1	GS01SB	GS02SB	GS03SB	GS04SB	GS05SB	GS06SB	GS07SB	GS08SB
% Moisture	%	18	19	28	23	19	26	26	17
4,4'-DDD (p,p'-DDD)	UG/KG	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
4,4'-DDE (p,p'-DDE)	UG/KG	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
4,4'-DDT (p,p'-DDT)	UG/KG	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
Aldrin .	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
alpha-BHC	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 ∪	2.3 U	2 U
alpha-Chlordane /2	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
beta-BHC	UG/KG	2.1 U	2.1 U	2.4 U	2.2 ∪	2.1 U	2.3 U ·	2.3 U	2 U
delta-BHC	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
Dieldrin	UG/KG	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
Endosulfan I (alpha)	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
Endosulfan II (beta)	UG/KG	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
Endosulfan Sulfate	UG/KG	4 U	4,1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
Endrin	UG/KG	4_Û	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 Ü
Endrin Aldehyde	UG/KG	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
Endrin Ketone	UG/KG	4 U	4.1 U	4.6 U	4.3 U	4.1 U	4.5 U	4.5 U	4 U
gamma-BHC (Lindane)	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
gamma-Chlordane /2	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
Heptachlor	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
Heptachlor Epoxide	UG/KG	2.1 U	2.1 U	2.4 U	2.2 U	2.1 U	2.3 U	2.3 U	2 U
Methoxychlor	UG/KG	21 U	21 U	24 U	22 U	21 U	23 U	23 U	20 U
PCB-1016 (Aroclor 1016)	UG/KG	40 U	41 U	46 U	43 U	41 U	45 U	45 U	40 U
PCB-1221 (Aroclor 1221)	UG/KG	82 U	83 U	93 U	87 U	83 U	91 U	91 U	81 U
PCB-1232 (Arocior 1232)	UG/KG	40 U	41 U	46 U	43 U	41 U	45 U	45 U	40 U
PCB-1242 (Aroclor 1242)	UG/KG	40 U	41 U	46 U	43 U	.41 U	45 U	45 U	40 U
PCB-1248 (Arocior 1248)	UG/KG	40 U	41 U	46 U	43 U	41 U	45 U	45 U	40 U
PCB-1254 (Aroclor 1254)	UG/KG	40 U	41 U	46 U	43 U	41 U	45 U	45 U	40 U
PCB-1260 (Aroclor 1260)	UG/KG	40 U	41 U	46 U	43 U	41 U	45 U	45 U	40 U
Toxaphene	UG/KG	210 U	210 U	240 U	220 U	210 U	230 U	230 U	200 U

	,	•								
Extractables Scan	. ,	GS01SB	GS02SE	377	GS03SB	GS04SB	GS05SB	GS06SB	GS07SB	GS08SB
% Moisture	%	18	19		28	23	19	26	26	17
3-and/or 4-)Methylphenol	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 (
1,1-Biphenyi	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 l
1,2,4,5-Tetrachiorobenzene	UG/KG	N	A	NA	NA	NA	NA NA	NA	. NA	
2,4,5-Trichlorophenol	UGAKG	1000 U	1000	טוו	1200 U	1100 U	1000 U	1100 U	1100 U	1000 (
2,4,6-Trichlorophenol	UG/KG	400 L		U	460 U	430 U	410 U	450 U	450 U	400 l
2,4-Dichlorophenol	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 (
2,4-Dimethylphenol	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400
2,4-Dinitrophenoi	UG/KG	1000 U	1000	UR	1200 U	1100 U	1000 U	1100 UJ	1100 UJ	1000 (
2,4-Dinitrotoluene	UG/KG	400 L	410	U	460 U	430 U	410 U	450 U	450 U	400 1
2,6-Dinitrotoluene	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U_	450 U	400 (
2-Chloronaphthalene	UG/KG	400 U		U	460 U	430 U	410 U	450 U	450 U	400
2-Chlorophenol	UG/KG	400 L	410	U	460 U	430 U	410 U	450 U	450 U	400 \
-Methyl-4,6-Dinitrophenol	UG/KG	1000 U	1000	U	1200 U	1100 U	1000 U	1100 UJ	1100 UJ	1000
?-Methylnaphthalene	UG/KG	400 L	410	U	460 U	430 U	410 U	450 U	450 U	400 (
?-Methylphenol	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 (
2-Nitroanlline	UG/KG	1000 U			1200 U	1100 U	1000 U	1100 U	1100 U	1000 (
2-Nitrophenol	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 l
3,3'-Dichlorobenzidine	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 1
-Nitroaniline	UG/KG	1000 U			1200 U	1100 U	1000 U	1100 U	1100 U	1000 (
-Bromophenyl Phenyl Ether	UG/KG	400 U	410	ilu l	460 U	430 U	410 U	450 U	450 U	400 L
-Chloro-3-Methylphenol	UG/KG	400 U	410	TO 1	460 U	430 U	410 U	450 U	450 U	400 (
-Chloroaniline	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 1
-Chlorophenyl Phenyl Ether	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 L
-Nitroaniline	UG/KG	1000 U	1000	UJ	1200 U	1100 U	1000 U	1100 U	1100 U	1000 1
Nitrophenol	UG/KG	1000 Ú	1000	UJ	1200 U	1100 U	1000 U	1100 U	1100 U	1000 L
cenaphthene	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
cenaphthylene	UG/KG	400 U	410	Ú	460 U	430 U	410 U	450 U	450 U	400 L
cetophenone	UG/KG	400 U	410	Ú	460 U	430 U	410 U	450 U	450 U	400 L
Anthracens	UG/KG	400 U	410	U	460 U	430 U	410 U.	450 U	450 U	74 J
Atrazine	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 L
Benzaldehyde	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 L
Benzo(a)Anthracene	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	120 J
Benzo(b)Fluoranthene	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	47 J	1000 J
Jenzo(ghi)Perylene	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	320 J
Benzo(k)Fluoranthene	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	45 J	980 J
Benzo-a-Pyrene	UG/KG	400 U	410	Ü	460 U	430 U	410 U	450 U	· 450 U	240 J
lenzyl Butyl Phthalate	UG/KG	400 U	J 410	Ü	460 UJ	430 UJ	410 UJ	450 U	450 U	400 L
is(2-Chloroethoxy)Methane	UG/KG	400 U	410	U	460 U	430 U	410 U	450 U	450 U	400 L
is(2-Chloroethyl) Ether	UG/KG	400 U	410	lu	460 U	430 U	410 U	450 U	450 U	400 L
is(2-Chloroisopropyl) Ether	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
Is(2-Ethylhexyl) Phthalate	UG/KG	400 U			460 UJ	430 UJ	450 J	450 U	450 U	400 L
aprolactam	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
arbazole	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U.	400 L
hrysene	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	220 J
ibenzo(a,h)Anthracene	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	150 J
ibenzofuran	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
iethyl Phthalate	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
imethyl Phthalate	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
i-n-Butyiphthalate	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
I-n-Octylphthalate	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
uoranthene	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	200 J
uorene	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 U
exachlorobenzene (HCB)	UG/KG	400 U	410		460 U	430 U	410 U	450 U	450 U	400 L
exachiorobenzene (nob)	UG/KG	400 U	410		460 U	430 U	410 U	450 U	450 U	400 U
exachlorocyclopentadiene (HCCP)	UG/KG	400 U			460 U	430 U	410 U	450 U	450 U	400 L
exachloroethane	UG/KG	400 U	410			430 U	410 U			
					460 U			450 U	450 U	400 L
deno (1,2,3-cd) Pyrene	UG/KG	400 U	410	(UJ	460 U	430 U	410 U	450 U	450 U	420

UG/KG

UG/KG

UG/KG

400 U

400 U 400 U

410 U

410 U

410 U

460 U 460 U 460 U

430 U 430 U 430 U

430 U

410 U

410 U

410 U

450 U

450 U

450 U

450 U

450 U

450 U

400 U 400 U

400 U

Isophorone Naphthalene

Nitrobenzene

Extractables Scan						•			
	<u> </u>	GS01SB	GS02SB	GS03SB	GS04SB	GS05SB	GS06SB	GS07SB	GS08SB
n-Nitrosodi-n-Propylamine	UG/KG	400 U	410 U	460 U	430 U	410 U	450 U	450 U	400 U
n-Nitrosodiphenylamine/Diphenylamine	UG/KG	400 U	410 U	460 U	430 U	410 U	450 U	450 U	400 U
Pentachlorophenol	UG/KG	1000 U	1000 UJ	1200 U	1100 U	1000 U	1100 U	1100 U	1000 U
Phenanthrene	UG/KG	400 U	410 U	460 U	430 U	410 U	450 U	450 Ü	400 U
Phenol	UG/KG	400 U	410 U	460 U	430 U	410 U	450 U	450 Ü	400 U
Pyrene	UG/KG	400 U	410 U	460 U	430 U	410 U	450 U	450 U	230 J
Volatiles Scan		GS01SB	GS02SB	GS03SB	GS04SB	GS05SB	GS06SB	GS07SB	GS08SB
% Moisture	%	18	19	28	23	19	. 26	26	17
1,1,1-Trichloroethane	UG/KG	11 0	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,1,2,2-Tetrachloroethane	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	UG/KG	11 U	11 U	23 Ü	15 U	12 U	13 U	13 U	10 U
1,1,2-Trichloroethane	UG/KG	11 0	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,1-Dichioroethane	UG/KG	11 0	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,1-Dichloroethene (1,1-Dichloroethylene)	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,2,3-Trichlorobenzene	UG/KG	NA	NA	NA	NA	NA NA	NA	NA	NA
1,2,4-Trichiorobenzene	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 Ú	10 U
1,2-Dibromo-3-Chloropropane (DBCP)	UG/KG	11 UJ	11 0	23 UJ	15 UJ	12 UJ	13 U	13 UJ	10 UJ
1,2-Dibromoethane (EDB)	UG/KG	11 0	11 0	23 U	15 U	12 U	13 U	13 U	10 U
1,2-Dichlorobenzene	UG/KG	11 0	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,2-Dichloroethane	UG/KG	11 0	11 0	23 U	15 U	12 U	13 U	13 U	10 U
1,2-Dichloropropane	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,3-Dichlorobenzene	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
1,4-Dichlorobenzene	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
Acetone	UG/KG	39 UJ	12 J	35 J	15 LU	12 UJ	21 J	13 UJ	99 J
Benzene	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
Bromochloromethane	UG/KG	NA NA	NA	NA	NA	. NA	NA	NA	NA
Bromodichloromethane	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
Bromoform	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
Bromomethane	UG/KG	11 UJ	11 UJ	23 UJ	15 UJ	12 UJ	13 U	13 UJ	10 UJ
Carbon Disuifide	.UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
Carbon Tetrachioride	UG/KG	11 U	11 U	23 U	15 U	12 U	13 U	13 U	10 U
Chlorobenzene	UG/KG	11 U	11 U	23 U	15 U	12 Ú	13 Ü	13 U	10 U
Chlorobenzene Chloroethane	UG/KG UG/KG	11 U	11 U	23 U 23 U	15 U 15 U	12 U 12 U	13 U 13 U	13 U 13 U	10 U
Chlorobenzene Chloroethane Chloroform	UG/KG	11 U	11 U	23 U	15 U	12 Ú	13 Ü	13 U	10 U
Chlorobenzene Chloroethane	UG/KG UG/KG	11 U 11 U 11 U	11 U 11 U 11 U	23 U 23 U 23 U	15 U 15 U 15 U	12 U 12 U 12 U	13 U 13 U 13 U	13 U 13 U 13 U	10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan	UG/KG UG/KG UG/KG	11 U 11 U 11 U GS015B	11 U 11 U 11 U	23 U 23 U 23 U GS03SB	15 U 15 U 15 U	12 U 12 U 12 U	13 U 13 U 13 U	13 U 13 U 13 U	10 U 10 U 10 U
Chlorobenzene Chloroform Volatlies Scan Chloromethane	UG/KG UG/KG UG/KG	11 U 11 U 11 U GS01SB	11 U 11 U 11 U 11 U GS02SB	23 U 23 U 23 U 23 U GS03SB 23 U	15 U 15 U 15 U 15 U GS04SB	12 U 12 U 12 U 12 U GS05SB	13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U	10 U 10 U 10 U GS08SB
Chlorobenzene Chloroform Volatiles Scan Chloromethane cis-1,2-Dichloroethene	UG/KG UG/KG UG/KG UG/KG	11 U 11 U 11 U 11 U GS015B	11 U 11 U 11 U 11 U GS02SB	23 U 23 U 23 U 23 U GS03SB 23 U 23 U	15 U 15 U 15 U 15 U GS04SB 15 U	12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U GS06SB 13 U 13 U	13 U 13 U 13 U 13 U GS07SB 13 U 13 U	10 U 10 U 10 U GS08SB 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene	UG/KG UG/KG UG/KG UG/KG	11 U 11 U 11 U 11 U GS015B 11 U 11 U	11 U 11 U 11 U 11 U GS02SB 11 U 11 U	23 U 23 U 23 U 23 U 33 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U GS0SSB 12 U 12 U 12 U	13 U 13 U 13 U 13 U GS06SB 13 U 13 U 13 U	13 U 13 U 13 U 13 U GS07SB 13 U 13 U	10 U 10 U 10 U GS08SB 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 33 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U GS06SB 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 6S07SB 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U GS08SB 10 U 10 U 10 U
Chlorobenzene Chlorochane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 33 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U GS06SB 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U GS07SB 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U GS08SB 10 U 10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 33 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U GS06SB 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 6S07SB 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U GS08SB 10 U 10 U 10 U
Chlorobenzene Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodiftuoromethane	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS01SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 33 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 6S06SB 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 6S07SB 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U GS08SB 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dichlorodifluoromethane Dichlorodifluoromethane Ethyl Benzene	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS01SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 33 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U GS06SB 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U .GS08SB 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodiftuoromethane Ethyl Benzene Isopropylbenzene	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS0158 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U GS08SB 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene Isopropylbenzene Methyl Acetate Methyl Butyl Ketone	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 16 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U 15	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 6S06SB 13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene Isopropylbenzene Methyl Acetate	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS0158 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene isopropyibenzene Methyl Acetate Methyl Butyl Ketone Methyl Ethyl Ketone Methyl Ethyl Ketone	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chloroethane Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifiuoromethane Ethyl Benzene Isopropylbenzene Methyl Acetate Methyl Ethyl Ketone Methyl Ethyl Ketone Methyl Isobutyl Ketone Methyl Isobutyl Ketone Methyl I-Butyl Ketone Methyl T-Butyl Ether (MTBE) Methylcyclohexane	UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropene Cyclohexane Dibromochloromethane Dichlorodifiuoromethane Ethyl Benzene Isopropylbenzene Methyl Acetate Methyl Butyl Ketone Methyl Sthyl Ketone Methyl Ishyl Ketone Methyl Ishyl Ketone Methyl Ishyl Ketone Methyl Ishyl Ketone	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chloroethane Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene Isopropylbenzene Methyl Acetate Methyl Ethyl Ketone Methyl Ethyl Ketone Methyl Isobutyl Ketone Methyl I-Butyl Ether (MTBE) Methylcyclohexane	UG/KG UG/KG	GS015B 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene Isopropylbenzene Methyl Acetate Methyl Acetate Methyl Sthyl Ketone Methyl Ishyl Ketone Methyl Isbyl Ketone Methyl Isbyl Ketone Methyl T-Butyl Ether (MTBE) Methylcyclohexane Methylene Chloride Styrene	UG/KG UG/KG	GS0158 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	11 U 11 U 11 U 11 U 11 U 11 U 11 U 11 U	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene isopropyibenzene Methyl Butyl Ketone Methyl Butyl Ketone Methyl I sobutyl Ketone Methyl I sobutyl Ketone Methyl T-Butyl Ether (MTBE) Methylcyclohexane Methylene Chloride Styrene Fetrachloroethene (Tetrachloroethylene)	UG/KG UG/KG	GS01SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 1	GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U
Chlorobenzene Chlorocthane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene Isopropylbenzene Methyl Acetate Methyl Butyl Ketone Methyl Ethyl Ketone Methyl Isobutyl Ketone Methylene Chloride Strene Tetrachloroethene (Tetrachloroethylene) Toluene	UG/KG UG/KG	GS015B   11 U   GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U	
Chloroethane Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene sopropylbenzene Methyl Benzene Methyl Butyl Ketone Methyl Ethyl Ketone Methyl Isobutyl Ketone Methyl I-Butyl Ether (MTBE) Methyloroethene	UG/KG UG/KG	GS015B   11 U   GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U	
Chlorobenzene Chloroethane Chloroform Volatiles Scan Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropene Cyclohexane Dibromochloromethane Dibromochloromethane Ethyl Benzene Isopropylbenzene Methyl Acetate Methyl Acetate Methyl Ketone Methyl Isbyl Ketone Methyl I-Butyl Ketone Methyl T-Butyl Ketone Methyl T-Butyl Ether (MTBE) Methylcyclohexane Methylene Chloride Styrene Tetrachloroethene (Tetrachloroethylene) Total Xylenes Trans-1,2-Dichloroethene	UG/KG UG/KG	GS015B   11 U   11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U	
Chloroethane Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene Isopropylbenzene Methyl Benzene Methyl Ethyl Ketone Methyl Ethyl Ketone Methyl Isobutyl Ketone Methylene Chloride Styrene Tetrachloroethene (Tetrachloroethylene) Total Xylenes Irans-1,2-Dichloropropene Irichloroethene (Trichloroethylene)	UG/KG UG/KG	GS01SB   11 U   GS02SB 11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U	
Chlorobenzene Chloroethane Chloroform Volatiles Scan  Chloromethane cis-1,2-Dichloroethene cis-1,3-Dichloropropene Cyclohexane Dibromochloromethane Dichlorodifluoromethane Ethyl Benzene isopropyibenzene Methyl Butyl Ketone Methyl Butyl Ketone Methyl Ethyl Ketone Methyl I sobutyl Ketone Methyl T-Butyl Ether (MTBE) Methylcyclohexane Methylene Chloride Styrene Tetrachloroethene (Tetrachloroethylene) Total Xylenes trans-1,2-Dichloroethene trans-1,3-Dichloropropene	UG/KG UG/KG	GS015B   11 U   11 U 11 U 11 U 11 U 11 U 11 U 11 U 11	23 U 23 U 23 U 23 U 23 U 23 U 23 U 23 U	15 U 15 U 15 U 15 U 15 U 15 U 15 U 15 U	12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	13 U 13 U 13 U 13 U 13 U 13 U 13 U 13 U	10 U 10 U 10 U 10 U 10 U 10 U 10 U 10 U	

Table 7. Subsurface Soll Analytical Results. Page 4 of 10

M	•	•	e	C	CB	n

		GS09SB	-	GS10SB		GS11SB		GS12SB		GS13SB	GS14SB	GS15SB	GS16SB	$\square$
% Moisture	%	20		18		18		17		15	15	21	23	
Aluminum	MG/KG	5800	5	8600	J	9600	_	3700	7	8400 J	4600 J	11000 J	12000	J
Antimony	MG/KG	0.6	U	0.96	R	1.6	R	0.58	כ	0.58 R	0.56 UJ	0.61 UJ		
Arsenic	MG/KG	0.88	C	13		6.6		4.6		4.7	7.4	8.1	7.3	
Barlum	MG/KG	29		430		41		19		23	54	45	49	
Beryllium	MG/KG	0.35		1.3		0.49		0.3		0.34	0.48	0.59	0.49	
Cadmium	MG/KG	0.05	U	0.87		0.05	U	0.05	اد	0.05 U	0.05 U	0.05 U	0.05	U
Calcium	MG/KG	540		9400		320		160		400	620	710	390	
Chromium	MG/KG	5.3		19		17		8.5		11	7.1	21	17	
Cobalt	MG/KG		R	12		3.2	R	1.4		1.7	12	2.9	2.1	
Copper	MG/KG	1.7	2	. 40	J	5.5	3	1.8		5.5 UJ	1.5 UJ	7 J	7.7	
Cyanide	MG/KG		NA		Ŋ		Š		NA	N/	NA NA	NA.		NA
Iron	MG/KG	3800		20000		20000		13000	,	14000	14000	24000	23000	
Lead	MG/KG	8.4		71		6.7		6.7		5.1	27	9.2	8.7	
Magnesium	MG/KG	290	7	2400	۲	680	J	170		470 J	250 J	640 J	650	
Manganese	MG/KG	9.3		1800		120		55		82	1000	410	52	
Nickel	MG/KG	2.6		20			٦	1.4	U	4.3 U	3.3 U	5.8 U	4.2	
Potassium	MG/KG	190		970	3	430		170		410 J	180 J	490 J	470	
Selenium	MG/KG	0.6		1.2		1.1		1.4		0.57 U	0.56 U	2.3 J	1.7	
Silver	MG/KG	0.18		0.76		1.1	R	0.47		0.39	0.42	0.65	0.45	
Sodium	MG/KG	440		490		350		96		310	260	290	290	
Thailium	MG/KG	0.78		0.76		0.75		0.75		0.73 U	0.73 U	0.79 U	0.8	
Total Mercury	MG/KG	0.06	UΠ	0.1	υΠ	0.1	U	0.06	U ¯	0.06 U	0.06 U	0.06 U	0.07	U
Vanadium Zinc	MG/KG MG/KG	8.4 12		30 290		33 44		26 8.4		22	23	36 36	33 25	

N/ Moleture		GS09SB	G\$10SB	GS11SB	GS12SB	GS13SB	GS14SB	G\$15SB	GS16SB
% Moisture	%	21	17	17	_26	15	15	21	22
4,4'-DDD (p,p'-DDD)	UG/KG	4.2 U	8.7 U	4 U	4.5 U	3.9 U	3.9 U	4.2 U	4.2 U
4,4'-DDE (p,p'-DDE)	UG/KG	4.2 U	4 U	1 4 U_	4.5 U	6.9	3.9 U	4.2 U	4.2 U
4,4'-DDT (p,p'-DDT)	UG/KG	4.2 U	110 N	4 U	4.5 U	3.9 U	16 J	4.2 U	4.2 U
Aldrin	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	2.2 U	2.2 U
alpha-BHC	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	2.2 U	2.2 ∪
alpha-Chlordane /2	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	2.2 U	2.2 U
beta-BHC	UG/KG	2.2 U	19 U	2 U	2.3 U	2 0	2 U	2.2 U	2.2 U
delta-BHC	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	2.2 U	2.2 U
Dieldrin	UG/KG	4.2 U	4 U	4 U	4.5 U	3.9 U	3.9 U	4.2 U	4.2 U
Endosulfan I (alpha)	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	2.2 U	2.2 U
Endosulfan II (beta)	UG/KG	4.2 U	4 U	4 U	4.5 U	3.9 U	3.9 U	4.2 U	4.2 U
Endosulfan Sulfate	UG/KG	4.2 U	4 U	4 U	4.5 U	3.9 U	3.9 U	4.2 U	4.2 U
Endrin	UG/KG	4.2 U	37 U	4 U	4.5 U	3.9 U	3.9 U	4.2 U	4.2 U
Endrin Aldehyde	UG/KG	42 U	15 U	4 U	4.5 U	3.9 U	3.9 U	4.2 U	4.2 U
Endrin Ketone	UG/KG	4.2 U	180 N	4 U	4.5 U	3.9 U	3.9 ∪	4.2 U	4.2 U
gamma-BHC (Lindane)	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	2.2 U	1.3 J
gamma-Chlordane /2	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	. 2.2 U	2.2 U
Heptachlor	UG/KG	2.2 U	_2 U	2 U	2.3 U	2 U	2 U	2.2 U_	2.2 U
Heptachlor Epoxide	UG/KG	2.2 U	2 U	2 U	2.3 U	2 U	2 U	2.2 U	2.2 U
Methoxychlor	UG/KG	22 Ü	360 U	20 U	23 U	20 U	43 U	22 U	22 U
PCB-1016 (Aroclor 1016)	UG/KG	42 U	40 U	40 U	45 U	39 U	39 U	42 U	42 U
PCB-1221 (Aroclor 1221)	UG/KG	85 U	81 U	81 U	91 U	79 U	79 U	85 U	86 U
PCB-1232 (Aroclor 1232)	UG/KG	42 U	40 U	40 U	45 U	39 U	39 U	42 U	42 U
PCB-1242 (Aroclor 1242)	UG/KG	42 U	40 U	40 U	45 U	39 U	39 U	42 U	42 U
PCB-1248 (Aroclor 1248)	UG/KG	42 U	40 U	40 U	45 U	39 U	39 U	42 U	42 U
PCB-1254 (Aroclor 1254)	UG/KG	42 U	40 U	40 U	45 U	39 U	. 39 U	42 U	42 U
PCB-1260 (Aroclor 1260)	UG/KG	42 U	40 U	40 U	45 U	39 U	39 U	42 U	70
Toxaphene	UG/KG	220 U	200 U	200 U	230 U	200 U	200 U	220 U	220 U

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	Scan

Extractables Scan	T	GS09SB	GS10SB	GS11SB	GS12SB	GS13SB	GS14SB	GS15SB	GS16SB
% Moisture	%	21	17	17	26	15	15	21	22
(3-and/or 4-)Methylphenol	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
1,1-Biphenyl	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
1,2,4,5-Tetrachlorobenzene	UG/KG	NA NA		NA.	NA	NA	NA	NA	NA
2,4,5-Trichlorophenol	UG/KG	1100 U	2000 U	1000 U	1100 U	980 U	980 U	1100 U	1100 U
2,4,6-Trichlorophenol	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
2,4-Dichlorophenol	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
2,4-Dimethylphenol	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
2,4-Dinitrophenol	UG/KG	1100 UJ	2000 UJ	1000 UJ	1100 UJ	980 U	980 UJ	1100 UR	1100 U
2,4-Dinitrotoluene	UG/KG	420 U	800 U	400 U	450 U	390 Ú	390 U	420 U	420 U
2,6-Dinitrotoluene	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U_
2-Chioronaphthalene	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
2-Chlorophenol	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
2-Methyl-4,6-Dinitrophenol	UG/KG	1100 U	2000 U	1000 U	1100 UJ	980 U	980 U	1100 U	1100 U
2-Methylnaphthalene	UG/KG	420 U	200 J	400 U	450 U	390 Ü	390 U	420 U	420 U
2-Methylphenol	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
2-Nitroaniline	UG/KG	1100 Ü	2000 U	1000 U	1100 U	980 U	980 U	1100 U	1100 U
2-Nitrophenol	UG/KG	420 U	800 U	400 U	. 450 U	390 U	390 U	420 U	420 U
3,3'-Dichlorobenzidine	UG/KG	420 U	800 U	400 U	450 U	390 UJ	390 U	420 U	420 UJ
3-Nitroaniline	UG/KG	1100 U 420 U	2000 U	1000 U	1100 U 450 U	980 U	980 U 390 U	1100 U	1100 U
4-Bromophenyl Phenyl Ether 4-Chloro-3-Methylphenol	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
4-Chloroaniline	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
4-Chlorophenyl Phenyl Ether	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
4-Nitroaniline	UG/KG	1100 U	2000 U	1000 U	1100 U	980 U	980 U	1100 UJ	1100 U
4-Nitrophenol	UG/KG	1100 U	2000 U	1000 U	1100 U	980 U	980 U	1100 UJ	1100 U
Acenaphthene	UG/KG	420 U	210 J	400 U	450 U	390 U	390 U	420 U	420 U
Acenaphthylene	UG/KG	420 U	6100	400 U	450 U	390 U	330 J	420 U	420 U
Acetophenone	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Anthracene	UG/KG	420 U	12000	49 J	450 U	390 U	690	420 U	420 U
Atrazine	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Benzaldehyde	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Benzo(a)Anthracene	UG/KG	420 U	23000	400 U	450 U	56 J	1400	420 U	420 U
Benzo(b)Fluoranthene	UG/KG	420 U	37000	400 U	450 U	160 J	2400	420 U	420 U
Benzo(ghi)Perylene	UG/KG	420 U	5200	400 U	450 U	390 U	540	420 U	420 U
Benzo(k)Fluoranthene	UG/KG	420 U	26000	46 J	450 U	170 J	2000	420 U	420 U
Benzo-a-Pyrene	UG/KG	420 U	25000	400 U	450 U	390 U	1400	420 U	420 U
Benzyl Butyl Phthalate	UG/KG	420 U 420 U	800 U	400 U	450 U 450 U	390 UJ	390 U	420 U	420 UJ 420 U
bis(2-Chloroethoxy)Methane	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
bis(2-Chloroethyl) Ether bis(2-Chloroisopropyl) Ether	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
bis(2-Ethylhexyl) Phthalate	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Caprolactam	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Carbazole	UG/KG	420 U	1800	400 U	450 U	390 U	210 J	420 U	420 U
Chrysene	UG/KG	420 U	35000	400 U	450 U	91 J	2100	420 U	420 U
Dibenzo(a,h)Anthracene	UG/KG	420 U	3300	400 U	450 U	390 U	390	420 UJ	420 U
Dibenzofuran	UG/KG	420 U	340 J	400 U	450 U	390 U	390 U	420 U	420 U
Diethyl Phthalate	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Dimethyl Phthalate	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Di-n-Butylphthalate	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Di-n-Octylphthalate	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Fluoranthene	UG/KG	420 U	28000	. 400 U	450 U	130 J	2200	420 U	420 U
Fluorene	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Hexachlorobenzene (HCB)	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 UJ	420 U
Hexachlorobutadiene	UG/KG	420 U	. 800 U	400 U	450 U	390 U	390 U	420 U	420 U
Hexachlorocyclopentadiene (HCCP)	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 UJ	420 U
Hexachloroethane	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Indeno (1,2,3-cd) Pyrene	UG/KG	420 U	20000	400 U	450 U	51 J	1100	420 UJ	420 U
Isophorone	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U
Naphthalene	UG/KG	420 U	390 J	400 U	450 U	390 U	390 U	420 U	420 U
Nitrobenzene	UG/KG	420 U	800 U	400 U	450 U	390 U	390 U	420 U	420 U

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X. Moleture         X. 23 (1988)         G\$1088											
n-Nicrosofth-Propylamine	Extractables Scan	<del></del>	GS00SB	Cetos	<u> </u>	G9119B	GS12SB	G9136B	GS14SB	CSTECH	COLCOR
Printenderghenylamina/Diphenylamine	n Alitenandi n Brandamina	LICING									
Pentenintence   USMG   420   U   2700   400   U   450   U   580   U   510   430   U   420   420   Mentenintence   USMG   420   U   420   Mentenintence   USMG   420   U   400   U   450   U   590   U   510   U   430   U   420   U   420   Mentenintence   USMG   420   U   400   U   450   U   590   U   500   U   400   U   420   U											
Penenthrene											
Phenol											
Votatiles Scan											
Volatiles Scan											
X. Moleture         X. 23 (1988)         G\$1088	Pyrene	UG/KG	420 (	3700	<u> </u>	40010	43010 1	89 7	24001	420 0	420 0
% Molature         %         21         17         17         26         15         16         21         22           1.1,12-Trichtoro-Starten         Gorgo         11         U 13         U 11         U 11         U 10         10         U 11         U 11         U 11         U 10         10         U 11	Voiatiles Scan	;	CCCCC		<u> </u>	004400	CC10CB	CC10CD	004400	·	COLCO
11,1-17-infororethane	B/ Malahum										
1,1,2,2-Fichero-Etane											
11,2-Trichicro-12,2-Triffuror-11    11,2-Trichicro-12,2-Triffuror-11    11,2-Trichicro-tethane											
11,12-Tichloroethane											
1.1-Dichloroethylene   USAG											
1.1-Dichloroethene (1.1-Dichloroethylene)											
12,3-Trichloroberszene											
12,4-Princhloropename (BGP)											11 U
1.2-Distromos-Schioropropane (DBCP)											N/
1,2-Dischorochenee											11 U
1,2-Dichlorobenzene											11 U.
1,2-Dichloropename											11 U
1,2-Dichioropropane											11 U
13-Dichlorobenzene											11 U
1.4-Dichloroberzene											11 U
Acetone											11 U
Benzene											11 U
Bromochloromethane											11 U.
Bromotichieromethane											11 U
Bromoform	Bromochloromethane										N/
Bromomethane	Bromodichloromethane										11 U
Carbon Tetrachloride	Bromoform										11 U
Carbon Tetrachloride	Bromomethane										<u>11   U.</u>
Chlorobenzene											11 U
Chloroethane											11 U
Chicroform											11 U
Chioromethane											11 U
Chioromethane		UG/KG	11   U	1	3 U_	11 U	11 U	10 U	10 U	11 U	11 U
Chloromethane	Volatiles Scan	ſ	GS09SB	65105	3	GS11SB	GS12SR	G\$135B	GS14SB	GS15SR	GS16SB
Cis-1,3-Dichloropropene	Chloromethane		11 U	1	3 U_	11 U	11 U	10 U	10 U	11 U	11 U
Cyclohexane											11 U
Dibromochloromethane											11 U
Dichlorodifiluoromethane			11 U	1	3 U			10 U	10 U	11 U	11 U
Ethyl Benzene											11 U
Isopropylbenzene											11 U
Methyl Acetate         Ug/KG         11 U         13 U         11 U         11 U         10 U         10 U         10 U         11 U											11 0
Methyl Butyl Ketone         UG/KG         11         UJ         13         U         11         UJ         11         UJ         10         U         11         U         11           Methyl Ethyl Ketone         UG/KG         11         UJ         13         U         11         UJ         10         U         11         U         11         UJ         11         UJ         10         U         11         U         11         UJ         10         U         11         UJ         11<	Isopropylbenzene	UG/KG	11 U	1	C	11 U	11 0	10 U	10 U	11 U	11 U
Methyl Ethyl Ketone         UG/KG         11 UJ         13 U         11 UJ         11 UJ         10 UJ         10 UJ         10 U         11 U         11 U         11 UJ         11 UJ         10 UJ         10 U         11 U         11 U         11 UJ         11 UJ         10 UJ         10 UJ         11 U         11 U         11 UJ         11 UJ         10 UJ         10 UJ         11 UJ         11 UJ         11 UJ         10 UJ         11 UJ         11 UJ         11 UJ         10 UJ         11 UJ	Methyl Acetate	UG/KG	11 0	1:	U	11 U	11 U	10 U	10 0	11 U	11 U
Methyl Isobutyl Ketone         UG/KG         11 UJ         13 U         11 UJ         11 UJ         10 UJ         10 UJ         10 UJ         11 UJ         11 UJ         11 UJ         10 UJ         10 UJ         11 UJ         11 UJ         11 UJ         10 UJ         11 UJ<	Methyl Butyl Ketone	UG/KG	11 U	J 1:	3 U	11 UJ	11 UJ	10 UJ	10 U	11 U	11 UJ
Methyl T-Butyl Ether (MTBE)         UG/KG         11 U         13 U         11 U         11 U         10 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         <	Methyl Ethyl Ketone	UG/KG	11 U	J 1:	I U	11 UJ	11 UJ	10 UJ	10 U	11 U	11 U.
Methyl T-Butyl Ether (MTBE)         UG/KG         11 U         13 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         <	Methyl Isobutyl Ketone	UG/KG									11 UJ
Methylcyclohexane         UG/KG         11 U         13 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U <th>Methyl T-Butyl Ether (MTBE)</th> <td>UG/KG</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>11 U</td>	Methyl T-Butyl Ether (MTBE)	UG/KG									11 U
Methylene Chloride         UG/KG         11 U         13 U         11 U         11 U         10 U         10 U         10 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U </td <th>Methylcyclohexane</th> <td>UG/KG</td> <td>11 0</td> <td></td> <td></td> <td>11 U</td> <td></td> <td></td> <td></td> <td></td> <td>11 U</td>	Methylcyclohexane	UG/KG	11 0			11 U					11 U
Styrene	Methylene Chloride										11 U
Tetrachloroethene (Tetrachloroethylene)	Styrene										11 U
Total Xylenes	Tetrachioroethene (Tetrachioroethylene)			1:	וייוו						11 U
Total Xylenes	Toluene										11 0
trans-1,2-Dichloroethene         UG/KG         11         U         13         U         11         U         11         U         10         U         10         U         11											11 0
trans-1,3-Dichloropropene         UG/KG         11 U         13 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         10 U         10 U         11 U         11 U         11 U         10 U         10 U         11 U <th< td=""><th>trans-1,2-Dichloroethene</th><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td>11 0</td></th<>	trans-1,2-Dichloroethene										11 0
Frichloroethene (Trichloroethylene)         UG/KG         11 U         13 U         11 U         11 U         10 U         10 U         11 U         11 U           Frichlorofluoromethane (Freen 11)         UG/KG         11 U         13 U         11 U         11 U         10 U         10 U         11 U         11 U	trans-1,3-Dichloropropene										11 0
Trichlorofluoromethane (Freon 11) UG/KG 11 U 13 U 11 U 11 U 10 U 10 U 11 U 11 U											11 0
											11 0
/Inyl Chloride UG/KG 11 U 13 U 11 U 10 U 10 U 11 U 11 U 11 U											

Table 7. Subsurface Soil Analytical Results. Page 7 of 10

		GS17SB	Γ	GS18SB		GS19SB		GS20SB		GS21SB	
% Moisture	%	21		18		20		16		20	
Aluminum	MG/KG	6700	J	7700	J	12000	5	4700	7	8600	J
Antimony	MG/KG	1.1	UJ	0.59	ŪJ	0.79	UJ	0.57	UJ	0.6	UJ
Arsenic	MG/KG	1.5		4.1		8.3		2.3		4.5	
Barlum	MG/KG	61		26	·	43		24		28	
Beryllium	MG/KG	0.39		0.31		0.46		0.3		0.3	
Cadmium	MG/KG	0.05	Ü	0.05	U	0.05	5	0.05	Ü		
Calcium	MG/KG	620		190		190		120		140	
Chromlum	MG/KG	7		8.7		12		4.7		14	
Cobalt	MG/KG	1.5		3.4		9.6		2.2		1.9	
Copper	MG/KG	2.4		3.9		6.3		2.3		3.6	
Cyanide	MG/KG		NA		NA		NA		NA		NA
Iron	MG/KG	5900		12000		20000		6900		13000	
Lead	MG/KG	9.1			J	12		6		6.8	
Magnesium	MG/KG	310	J	470	J	680	J	210	J	420	J
Manganese	MG/KG	57		76		600		140		95	
Nickel	MG/KG	2.5		3.4		5.4		1.7		3.6	
Potassium	MG/KG	280		220		480		170		300	
Selenium	MG/KG	0.6		0.59		1.7	J	0.57		0.95	R
Silver	MG/KG	0.32		0.32	R	0.53		0.25	R	0.4	
Sodium	MG/KG	340		540		280		340		310	
Thallium	MG/KG	0.78		0.76		0.78		0.74		0.78	
Total Mercury	MG/KG	0.06	U	0.06	٦	0.06	U_	0.06	U	0.06	U
Vanadlum	MG/KG	13		18		31		14		21	
Zinc	MG/KG	12		16		30		11		16	

	L .	GS17SB		GS18SB		GS19SB		GS20SB		GS21SB
% Moisture	%	22		25		21		18		20
4,4'-DDD (p,p'-DDD)	UG/KG	4.2	Ü	4.4	U	4.2	Ū	4	ς	4.1 U
4,4'-DDE (p,p'-DDE)	UG/KG	4.2	U	4.4	U	4.2		4	Ü	4.1 U
4,4'-DDT (p,p'-DDT)	UG/KG	4.2	Ü	4.4	U.	4.2	ū	4	C	4.1 U
Aldrin	UG/KG	2.2	U	2.3	U	2.2	Ü	2.1	C	2.1 U
alpha-BHC	UG/KG	2.2	U	2.3	U	2.2	Ú	2.1	U	2.1 U
alpha-Chlordane /2	UG/KG	2.2	υ	2.3	U	2.2	Ū	2.1	Ū	2.1 U
beta-BHC	UG/KG	2.2	U	2.3	U	2.2	Û	2.1	Ū	2.1 U
delta-BHC	UG/KG	2.2	U	2.3	U	2.2	Ū	2.1	Ū J	2.1 U
Dieldrin	UG/KG	4.2	U	4.4	U	4.2		4	U	4.1 U
Endosulfan I (alpha)	UG/KG	2.2	U	2.3	U	2.2	Ū	2.1	Ū	2.1 U
Endosulfan II (beta)	UG/KG	4.2	U	- 4.4	Ū	4.2	Ū	4	U	4.1 U
Endosulfan Sulfate	UG/KG	4.2	υ	4.4	U	4.2	Ū	4 1	Ū	4.1 U
Endrin	UG/KG	4.2	U	4.4	U	4.2	Ū	4 (	U	4.1 U
Endrin Aldehyde	UG/KG	4.2	Ū	4.4	U	4.2	Ū	4 1	U	4.1 U
Endrin Ketone	UG/KG	4.2	U	4.4	Ū	4.2	Ū	4 (	U	4.1 U
gamma-BHC (Lindane)	UG/KG	2.2	U	1.7	J	2.2	Ŭ	2.1	Ū	2.1 U
gamma-Chlordane /2	UG/KG	2.2	c	2.3	U	2.2	Ū	2.1	Ü	2.1 U
Heptachlor	UG/KG	2.2	c	2.3	U	2.2	Ū	2.1	J	2.1 U
Heptachlor Epoxide	UG/KG	2.2	C	2.3	U	2.2	Ū	2.1	J	2.1 U
Methoxychlor	UG/KG	22	U	23	Ū	22	Ū	21	J	21 U
PCB-1016 (Arocior 1016)	U\$/KG	42	U	44	U	42	Ū	40 1	J	41 U
PCB-1221 (Araclor 1221)	UG/KG	86	U	89	c	85	J	82 (	j -	84 U
PCB-1232 (Aroclor 1232)	UG/KG	42	Ū	44	حا	42	Ū	40 (	7	41 U
PCB-1242 (Aroclor 1242)	U6/KG	42	Ū	44	Ū	42	J	40 L	1	41 U
PCB-1248 (Aroclor 1248)	UG/KG	42	Ū	44	Ü	42	J	40 L	7	41 U
PCB-1254 (Aroclor 1254)	UG/KG	42	U	44	Ū	42	Ĵ	40 L	1	41 U
PCB-1250 (Aroclor 1260)	UG/KG	42	Ū	42	J	39 .	J	40 (	1	41 U
Toxaphene	UG/KG	220	Ū	230	Ū	220	<u>,</u>	210 (	7	210 U

Extractables Scan

Extractables Scan											
ļ.,		GS17SB		GS18SB		G\$19SB		GS20SB		GS21SB	-
% Molsture	%	22		25		21	<del> </del>	18		20	
(3-and/or 4-)Methylphenol	UG/KG	420		440		420		400		410 L	
1,1-Biphenyl	UG/KG	420		440		420		400		410 U	
1,2,4,5-Tetrachiorobenzene	UG/KG		NA		NA	<del> </del>	NA	4000	NA		١A
2,4,5-Trichlorophenol	UG/KG	1100			U	1100		1000		1000 U	
2,4,6-Trichlorophenol	UG/KG	420		440		420	U	400		410 L	_
2,4-Dichlorophenol	UG/KG	420	5	440	U	420		400		410 U	
2,4-Dimethylphenol	UG/KG	420		440		420		400		410 L	
2,4-Dinitrophenoi	UG/KG	1100		1100		1100		1000		1000 L	
2,4-Dinitrotoluene	UG/KG	420		440		420		400		410 0	
2,6-Dinitrotoluene	UG/KG	420		440		420		400		410 U	
2-Chloronaphthalene	UG/KG	420		440		420		400		410 L	
2-Chlorophenol	UG/KG	420		440		420		400 1000		410 U	
2-Methyl-4,6-Dinitrophenol	UG/KG	1100		1100		1100					
2-Methylnaphthalena	UG/KG	420		440		420		400		410 U	
2-Methylphenol 2-Nitroaniline	UG/KG	420 1100		440 1100		420 1100		400 1000		1000 U	
2-Nitrophenol	UG/KG UG/KG	420		440		1100		400		410 L	
3,3'-Dichlorobenzidine	UG/KG	420		440		420		400		410 0	
3-Nitroaniline	UG/KG	1100		1100		1100		1000		1000 U	
4-Bromophenyl Phenyl Ether	UG/KG		Ü	440		420		400		410 U	
4-Chloro-3-Methylphenol	UG/KG	420		440		420		400		410 U	
4-Chloroaniline	UG/KG	420		440		420		400		410 U	
4-Chlorophenyl Phenyl Ether	UG/KG	420		440		420		400		410 U	
4-Nitroaniline	UG/KG	1100		1100		1100		1000		1000 U	
4-Nitrophenol	UG/KG	1100		1100		1100		1000		1000 U	
Acenaphthene	UG/KG	420		440		420		400		410 U	
Acenaphthylene	UG/KG	420			ŭ	420		400		410 U	
Acetophenone	UG/KG	420		440	Ü	420		400		410 U	$\sqcap$
Anthracene	UG/KG	420	U	440	Ü	420	5	400	υT	410 U	,
Atrazine	UG/KG	420	U	440	J	420	5	400	U	410 U	ī
Benzaldehyde	UG/KG	420	C	440	C	420	c	400	U	410 U	亓
Benzo(a)Anthracene	UG/KG	420	U	440	J	420	U	400	U	410 U	$\overline{I}$
Benzo(b)Fluoranthene	UG/KG		J	440	٥	420		400		410 U	
Benzo(ghi)Perylene	UG/KG	420			٦	420		120		410 U	
Benzo(k)Fluoranthene	UG/KG		J		U	420		400		410 U	
Benzo-a-Pyrene	UG/KG	420		440	٦	420	υ	400		410 U	
Benzyl Butyl Phthalate	UG/KG	420			UJ	420		400		410 U	
bis(2-Chloroethoxy)Methane	UG/KG	420			U	420		400		410 U	
bis(2-Chloroethyi) Ether	UG/KG	420			U	420		400		410 U	
bis(2-Chlorolsopropyl) Ether	UG/KG	420			U	420		400		410 U	
bis(2-Ethylhexyl) Phthalate	UG/KG	420			IJ	420			Ü	410 U	
Caprolactam Carbazole	UG/KG	420			U	420		400		410 U	
Chrysene	UG/KG	420 420			U	420 420	U	400 400		410 U	
Dibenzo(a,h)Anthracene	UG/KG	420			Ü	420 420		400		410 U	
Dibenzofuran	UG/KG	420			ö		U	400		410 U	
Diethyl Phthalate	UG/KG	420		440			U	400		410 U	
Dimethyl Phthalate	UQ/KG	420		440		420		400		410 U	
Di-n-Butylphthalate	UG/KG	420		440			<del>ö</del>	400		410 U	
Di-n-Octylphthalate	UG/KG	420		440		420			<del>ŭ</del> †	410 U	
Fluoranthene	UG/KG	420		440			υ	400		410 U	
Fluorene	UG/KG	420		440			υ	400		410 U	
Hexachlorobenzene (HCB)	UG/KG	420		440		420		- 400		410 U	
Hexachlorobutadiene	UG/KG	420		440		420		400		410 U	
Hexachlorocyclopentadiene (HCCP)	UG/KG	420		440		420		400		410 U	
Hexachloroethane	UG/KG		Ū	440		420		400		410 U	
Indeno (1,2,3-cd) Pyrene	UG/KG	420		440		420		400		410 U	
Isophorone	UG/KG	420		440		420		400		410 U	
Naphthalene	UG/KG	420	υİ	440	υl	420		400		410 U	
Nitrobenzene	UG/KG	420	υT	440	U	420		400		410 U	$\neg$
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		GS17SB	7	GS18SB		GS19SB		GS20SB	GS21SB
n-Nitrosodi-n-Propylamine	UG/KG	420 U	T	440	U	420	U	400 U	410 U
n-Nitrosodiphenylamine/Diphenylamine	UG/KG	420 U	T	440	C	420	U	400 U	410 U
Pentachlorophenol	UG/KG	1100 U		1100	U	1100	U	1000 U	1000 U
Phenanthrene	UG/KG	420 U	ıT	440	U	420	Ū_	400 U	410 U
Phenol	UG/KG	420 U		440	U_	420	U	400 U	410 U
Pyrene	UG/KG	420 U		440	U	420	U	400 U	410 U

# Volatiles Scan

,	. Г	GS17SB	_	GS18SB	Γ	GS19SB		GS20SB	GS21SB
% Moisture	%	22		25		21	_	18	20
1,1,1-Trichioroethane	UG/KG	13	U	11	U	11	U	11 U	11 U
1,1,2,2-Tetrachloroethane	UG/KG	13	U	11	U	11	U	11 U	11 U
1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	UG/KG	13	Ū	11	U	11	U	11 U	11 U
1,1,2-Trichloroethane	UG/KG	13	U	11	U	11	U	11 U	11 U
1,1-Dichloroethane	UG/KG	13	U	11	U	11	U	11 U	11 U
1,1-Dichloroethene (1,1-Dichloroethylene)	UG/KG	13	U	11	U	. 11	U	11 U	11 U
1,2,3-Trichiorobenzene	UG/KG		NA		NA		NA	N/	NA NA
1,2,4-Trichiorobenzene	UG/KG	13	U	11	دا	11	Ü	11 U	11 U
1,2-Dibromo-3-Chloropropane (DBCP)	UG/KG	13	U	11	IJ	11	IJ	11 U	11 U
1,2-Dibromoethane (EDB)	UG/KG	13	U	11	c	11	C	11 U	11 U
1,2-Dichlorobenzene	UG/KG	13	U	11	U_	11	Ü	11]U	_11 U
1,2-Dichloroethane	UG/KG	13	U	11	U	11	C	11 U	11 U
1,2-Dichloropropane	UG/KG	13	U	11	c	11	Ü	11 U	11 U
1,3-Dichlorobenzene	UG/KG	13		11	U	.11	U	11 U	11 U
1,4-Dichlorobenzene	UG/KG	13	٥	11	٦	11	U	11 U	11 U
Acetone	UG/KG	13		. 14	IJ	26	IJ	11 U	11 UJ
Benzene	UG/KG	13		11	υ	11		11 U	11 U
Bromochloromethane	UG/KG		NA		NA		NΑ	N/	
Bromodichioromethane	UG/KG	13		11		11		11 U	. 11 U
Bromoform	UG/KG	13			U	11	U	11 U	11 U
Bromomethane	UG/KG		3		UJ		3	11 U.	
Carbon Disulfide	UG/KG	13		11		11		11 U	11 U
Carbon Tetrachloride	UG/KG	13		11		11		11 U	11 U
Chlorobenzene	UG/KG	13			U			11 U	11 U
Chloroethane	UG/KG				U	11		11 U	11 U
Chloroform	UG/KG	13	U	11	U	11	U	11 0	11 0

# Volatiles Scan

	[	GS17SB	GS18SB	GS19SB	GS20SB	GS21SB
Chloromethane	UG/KG	13 U	11 U.	11 U	11 U	11 U
cis-1,2-Dichloroethene	UG/KG	13 U	11 U	11 U	11 0	11 U
cis-1,3-Dichioropropene	UG/KG	13 U	11 U	11 U	11 U	11 U
Cyclohexane	UG/KG	13 U	11 U	11 U	11 U	11 U
Dibromochioromethane	UG/KG	13 U	11 U	11 U	11 0	11 U
Dichlorodifluoromethane	UG/KG	13 U	11 U	11 U	11 U	11 U
Ethyl Benzene	UG/KG	13 U	· 11 U	11 U	11 U	11 U
Isopropyibenzene	UG/KG	13 U	11 U	11 U	11 U	11 U
Methyl Acetate	UG/KG	13 U	11 U	11 U	11 U	11 0
Methyl Butyl Ketone	UG/KG	13 U	11 UJ	11 UJ	11 U	11 UJ
Methyl Ethyl Ketone	UG/KG	13 U	11 UJ	11 UJ	11 U	11 UJ
Methyl Isobutyl Ketone	UG/KG	13 U	11 UJ	11 UJ	11 U	11 00
Methyl T-Butyl Ether (MTBE)	UG/KG	13 U	11 U	11 U	11 U	11 U
Methylcyclohexane	UG/KG	13 U	11 U	11 0	11 U	11   U
Methylene Chloride	UG/KG	13 U	11 U	1.1 U	11 U	11 U
Styrene	UG/KG	13 U	11 U	11 U	11 U	11 U
Tetrachloroethene (Tetrachloroethylene)	UC/KG	13 U	11 U	11 U	11 U	11 U
Toluene	UG/KG	13 U	11 U	11 U	11 U	11 U
Total Xylenes	UG/KG	13 U	11 U	11 U	11 U	11 U
trans-1,2-Dichloroethene	UG/KG	13 U	11 0	11 U	11 U	11 U
trans-1,3-Dichioropropene	UG/KG	13 U	11 U	11 0	11 U	11 U
Trichioroethene (Trichioroethylene)	UG/KG	13 U	11 U	11 0	11 U	11 U
Trichlorofluoromethane (Freon 11)	UG/KG	13 U	11 U	11 U	11 U	11 U
Vinyl Chloride	UG/KG	13 U	11 U	11 U	11 U	11 U

#### **Data Qualifiers**

U-Analyte not detected at or above reporting limit. The number is the minimum quantitation limit.

J-identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-identification of analyte is acceptable; reported value may be blased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be blased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolities of technical chlordane

Table 8. Sediment Analytical Results. Page 1 of 5

## Metals Scan

·		GS01SD		GS02SD		GS03SD		GS04SD		GS05SD		GS06SD		GS07SD		GS08SD	
% Moisture	%	61		56		84		77		54		20		22		27	
Aluminum	MG/KG	8200	<u>_</u>	4700	J	16000	J	9100	J	4900	J	450	J	140 J		510	J
Antimony	MG/KG	1.2	UJ	2.4	UJ	4.5	3	2.1	UJ	2.1	R_	0.7	R_	0.61 L		0.65	
Arsenic	MG/KG	5.7		1.8	R	7.9	R	5.5		2.9	R	0.88	U_	0.9 U		0.95	U
Barlum	MG/KG	89		77		300		120		71		6.8		3.2		15	
Beryllium	MG/KG	0.39		0.56		1.3		1.2		0.51		0.03	U	0.03 U		0.1	1
Cadmium	MG/KG	0.1		0.09	Ū	0.25	כ	0.18	U	0.09	U_	0.05	U_	0.05 U		0.06	U
Calcium	MG/KG	1500		880		2000		2000		1200		250		110 L		180	
Chromium	MG/KG	13		7.8		24		15		8.4		1.3		0.61 U		1.4	
Cobalt	MG/KG	4.4		4.7		12		11		5.5		0.61	R	0.47		1	
Copper	MG/KG	10	Ū	6.6	U_	. 25	U	12	UJ	5.8	UJ	0.15	S	0.22 U	IJ	1.2	
Cyanide	MG/KG		NA		NA		NA		NA		NA		NA	N	IÄ		NA
Iron	MG/KG	15000		6400		24000		13000		8400		950		340		2200	J
Lead	MG/KG	12	J	20	J	41	J	21		12		1.8		1.2 J		3.6	
Magneslum	MG/KG	1000	J	410	J	1200	J	1000		660	J	54	J	14 U	IJ	81	
Manganese	MG/KG	340		500		720		280		320		52		53		44	
Nickel	MG/KG	6.6		4.6		16		15		6.5		0.76		0.6 U		1	U
Potassium	MG/KG	700		410		1100	J	811		480		90		52 J		56	
Selenium	MG/KG	1.9		1.1		4.2		2.1			U	0.6		0.61 U		0.65	
Silver	MG/KG	0.68	R	0.45	R	1.1		1.1		0.71	R	0.24	R_	0.18 U		0.19	Ü
Sodium	MG/KG	690		580		1500		1200		600		330		330		68	
Thailium	MG/KG	1.6		1.4		3.9	U	2.8		1.3		0.78		0.79 U		0.85	
Total Mercury	MG/KG	0.13	U.	0.11	U	0.31	U	0.22		0.11	U_	0.06		0.06 U		0.07	<u>u</u>
Vanadium	MG/KG	24		13	-	41		23		13		1.6	R	0.57 R		2.4	
Zinc	MG/KG	36		33		130		84	l	39		3.6		2.9 U		4.6	

#### Pesticides & Aroclors Scan

		GS01SD	GS02SD	GS03SD	GS04SD	GS05SD	GS06SD	GS07SD	GS08SD
% Moisture	%	55	52	79	78	54	20	18	19
4,4'-DDD (p,p'-DDD)	UG/KG	7.3 U	6.9 U	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
4,4'-DDE (p,p'-DDE)	UG/KG	7.3 U	7.9	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
4,4'-DDT (p,p'-DDT)	UG/KG	7.3 U	6.9 U	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
Aidrin	UG/KG	3.8 ∪	3.5 U	8.1 U	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
alpha-BHC	UG/KG	3.8 U	3.5 U	8.1 U	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
alpha-Chlordane /2	UG/KG	3.8 U	3.5 U	8.1 U	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
beta-BHC	UG/KG	3.8 U	3.5 U	12	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
delta-BHC	UG/KG	3.8 U	3.5 U	8.1 U	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
Dieldrin	UG/KG	7.3 U	6.9 U	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
Endosulfan I (alpha)	UG/KG	3.8 U	3.5 U	8.1 Ü	7.7 U	3.7 U	2.1 U .	2.1 U	2.1 U
Endosulfan II (beta)	UG/KG	7.3 U	6.5 J	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
Endosulfan Sulfate	UG/KG	7.3 U	10 N	14 NJ	15 U	7.2 U	4.1 U	4 U	4.1 U
Endrin	UG/KG	7.3 U	6.9 U	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
Endrin Aldehyde	UG/KG	7.3 U	6.9	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
Endrin Ketone	UG/KG	7.3 U	6.9 U	16 U	15 U	7.2 U	4.1 U	4 U	4.1 U
gamma-BHC (Lindane)	UG/KG	3.8 U	3.5 U	8.1 U	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
gamma-Chlordane /2	UG/KG	3.8 U	3.5 U	8.1 U	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
Heptachlor	UG/KG	3.8 U	3.5 U	8.1 U	7.7 U	3.7 U	2.1 U	2.1 U	2.1 U
Heptachlor Epoxide	UG/kG	3.8 U	3.5 U	8.1 U	7.7 U	3.7 ∪	2.1 U	2.1 U	2.1 Ú
Methoxychlor	UG/KG	38 U	35 U	81 U	77 U	37 U	21 U	21 U	21 U
PCB-1016 (Aroclor 1016)	UG/KG	73 U	69 U	160 U	150 U	72 U	41 U	40 U	41 U
PCB-1221 (Aroclor 1221)	UG/KG	150 U	140 U	320 U	300 U	150 U	84 U	82 U	83 U
PCB-1232 (Aroclor 1232)	UG/KG	73 U	. 69 U	160 U	150 U	72 U	41 U	40 U	41 U
PCB-1242 (Aroclor 1242)	UG/KG	73 U	69 U	160 U	150 U	72 U	41 U	40 U	41 U
PCB-1248 (Aroclor 1248)	UG/KG	73 U	69 U	160 U	150 U	72 U	41 U	40 U	41 U
PCB-1254 (Aroclor 1254)	UG/KG	73 U	69 U	160 U	150 U	72 U	41 U	40 U	41 U
PCB-1260 (Aroclor 1260)	UG/KG	73 U	69 U	160 U	150 U	72 U	41 U	40 U	41 U
Toxaphene	UG/KG	380 U	350 U	810 U	770 U	370 U	210 U	210 U	210 U

## Extractables Scan

	:	GS01SD	GS02SD	GS03SD	GS04SD	GS05SD	GS06SD	GS07SD	GS08SD
% Moisture	%	55	52	79	78	54	20	18	19
(3-and/or 4-)Methylphenol	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
1,1-Biphenyl	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
1,2,4,5-Tetrachiorobenzene	UG/KG	NA_	NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA
2,4,5-Trichlorophenol	UG/KG	1100 U	1000 U	2400 U	2300 U	1100 U	830 U	830 U	830 U
2,4,6-Trichlorophenol	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 Ù	330 U
2,4-Dichlorophenol	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
2,4-Dimethylphenol	UG/KB	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
2,4-Dinitrophenol	UG/KG	1100 U	1000 U	2400 U	2300 U	1100 UR	830 UJ	830 UR	830 UJ
2,4-Dinitrotoluene	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
2,6-Dinitrotoluene	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
2-Chloronaphthalene	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
2-Chlorophenol	UG/KG	440 U	410 U	940 U 2400 U	900 U 2300 U	430 U 1100 U	330 U	330 U 830 U	330 U 830 U
2-Methyl-4,6-Dinitrophenol	UG/KG UG/KG	1100 U	1000 U 410 U	930 U	900 U	430 U	830 U	330 U	330 U
2-Methylnaphthalene 2-Methylphenol	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
2-Nitroaniline	UG/KG	1100 U	1000 U	2400 U	2300 U	1100 U	830 U	830 U	830 U
2-Nitrophenol	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
3,3'-Dichiorobenzidine	UG/KG	440 U	410 UJ	940 UJ	900 UJ	430 U	330 U	330 U	330 U
3-Nitroaniline	UG/KG	1100 U	1000 U	2400 U	2300 U	1100 U	830 U	830 U	830 U
4-Bromophenyl Phenyl Ether	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
4-Chloro-3-Methylphenol	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
4-Chioroaniline	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
4-Chlorophenyl Phenyl Ether	UG/KG	440 U	410 U	940 U	900 Ü	430 U	330 U	330 U	330 U
4-Nitroaniline	UG/KG	1100 U	1000 U	2400 U	2300 U	1100 UJ	830 U	830 UJ	830 U
4-Nitrophenol	UG/KG	1100 U	1000 U	2400 U	2300 U	1100 UJ	830 U	830 UJ	830 U
Acenaphthene	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Acenaphthylene	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Acetophenone	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Anthracene	UG/KG	. 440 U	410 U	270 J	900 U	430 U	330 U	330 U	330 U
Atrazine	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Benzaldehyde	UG/KG	89 J	410 U	110 J	130 J	430 U	330 U	330 U	330 U
Benzo(a)Anthracene	UG/KG	440 U	410 U	150 J	900 U	430 U	330 U	330 U	330 U
Benzo(b)Fluoranthene	UG/KG	440 U	410 U	540 J	900 U	430 U	330 U	330 U	330 U
Benzo(ghi)Perylene	UG/KG	92 J	410 U	940 U	900 U	430 U	330 U	330 U	330 U.
Benzo(k)Fluoranthene	UG/KG	440 U	410 U	560 J	900 U	430 U	330 U	330 U	330 U
Benzo-a-Pyrene	UG/KG	440 U	410 U 410 UJ	150 J	900 UJ	430 U	330 U	330 U 330 U	330 U
Benzyl Butyl Phthalate bis(2-Chloroethoxy)Methane	UG/KG UG/KG	440 UJ 440 U	410 U	940 UJ 940 U	900 U	430 U	330 U	330 U	330 U
bis(2-Chloroethyl) Ether	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
bis(2-Chloroisopropyl) Ether	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
bis(2-Ethylhexyl) Phthalate	UG/KG	440 UJ	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Caprolactam	UG/KG	440 U	410 U	940 Ú	900 U	430 U	330 U	330 U	330 U
Carbazole	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Chrysene	UG/KG	440 U	410 U	250 J	900 U	430 U	330 U	330 U	330 U
Dibenzo(a,h)Anthracene	UG/KG	440 U	410 U	940 U	900 U	430 UJ	330 Ü	330 UJ	330 U
Dibenzofuran	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Diethyl Phthalate	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Dimethyl Phthalate	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Di-n-Butylphthalate	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Di-n-Octylphthalate	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Fluoranthene	UG/KG	440 U	410 U	340 J	900 U	430 U	330 U	330 U	330 U
Fluorene	UG/KG	440 U	410 U	130 U	900 U	430 U	330 U	330 U	330 U
Hexachiorobenzene (HCB)	UG/KG	440 U	410 U	940 U	900 U	430 UJ	330 U	330 UJ	330 U
Hexachlorobutadiene	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Hexachlorocyclopentadiene (HCCP)	UG/KG	440 U	410 U	940 U	900 U	430 UJ	330 U	330 UJ	330 U
Hexachloroethane	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Indeno (1,2,3-cd) Pyrene	UG/KG	440 U	410 U	180 J	900 U	430 UJ	330 U	330 UJ	330 U
sophorone	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Naphthalene	UG/KG	440 U	· 410 U	940 U	3 J	430 U	330 U	330 U	330 U
Nitrobenzene	UGIKG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
n-Nitrosodi-n-Propylamine	UG/KG	440 U	410 U	940 Ū	900 U	430 U	330 U	330 U	330 U
n-Nitrosodiphenylamine/Diphenylamine	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U

Extractables Sca	ſ	ľ																																									ĺ					ı			ı	ı			ı	ı	ı	ı	ı	ı	ı																			ı	ı	ı										ı	ı	ı												١	١			ļ	ı				į													١				ĺ								
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		GS01SD	GS02SD	GS03SD	GS04SD	GS05SD	GS06SD	GS07SD	GS08SD
Pentachlorophenol	UG/KG	_1100 U	1000 U	190 U	2300 U	1100 UJ	25 U	830 UJ	830 U
Phenanthrene	UG/KG	440 U	410 U	120 J	900 U	430 U	330 U	330 U	330 U
Phenol	UG/KG	440 U	410 U	940 U	900 U	430 U	330 U	330 U	330 U
Pyrene	UG/KG	440 U	410 U	220 J	120 J	430 U	330 U	330 U	330 U

Misodillitoda Extragasica	- 1	GS01SD		GS02SD	-	GS03SD		GS04SD		GS05SD		GS06SD	GS075	D	GS08SD
.GAMMASITOSTEROL	UG/KG		NR	1400	NJ		NR		NR		NR	N	3	NR_	NR
17 UNKNOWN COMPOUNDS	UG/KG		NR		NA I		NR		NR		NR	N	1	NR	NR
1-DOCOSENE	UG/KG		NR		NR		NR		NR		NR	N N	₹	NR	NR
1-HEXADECENE	UG/KG	540	NJ		NR	1000	NJ		NR	1	NR	N.		NR	NR
23 UNKNOWN COMPOUNDS	UG/KG	23000	J		NR		NR		NR		NR	N		NR	NR
24 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR	54000	J		NR	N	1	NR	NR
25 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR		NR	140000	J	N		NR	NR
25 UNKNOWNS	UG/KG		NR		NR	54000	J		NR		NR_	N N	}	NR	NR.
28 UNIDENTIFIED COMPOUNDS	UG/KG		NR	92000	J		NR		NR		NR	N	1	NR	NR
3 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR		NR		NR	N	?	NR	170 J
BENZENE, 1-METHYL-2-ISOPROPYL	UG/KG		NR		NR		NR		NR_	1200	NJ	N	1	NR	NR
BENZO [E] PYRENE	UG/KG		NR		NR	590	NJ		NR		NR	N	₹	NR	NR
CHOLESTANOL	UG/KG		NR		NR		NR	1600	ÑJ		NR	N	1	NR	NR.
ERGOST-5-EN-3-OL, (3.BETA.)-	UG/KG		NR		NR		NR	1700	NJ		NR_	N		NR	NR
ERGOST-5-N-3-OL, (3.BETA.)-	UG/KG	1500			NR		NR		NR		NR	N		NR	NR
HEXADECANOIC ACID	UG/KG	420	NJ		NR :		NR		NR		NR	N		NR	NR
PHENANTHRENONE DERIVATIVE	UG/KG		NR		NR:		NR		NR	2500		N		NR	NR
STIGMAST-4-EN-3-ONE	UG/KG	730			NR	1400	NJ_	2200	NJ	800	NJ_	N		NR	NR.
STIGMASTEROL	UG/KG	1100			NR		NR		NR		NR	N.		NR	NR
UNKNOWN ALCOHOL	UGAKG		NR		NR		NR	9200			NR_	N		NR	NA
UNKNOWN CARBOXYLIC ACID	UG/KG		NR		NE		NR		NR	500	J	N		NR	NR
UNKNOWN COMPOUND	UG/KG		NR		NR		NR		NR		NR	61 J		9 )	NR
UNKNOWN KETONE	UG/KG		NR		NR		NA	1800	7		NR	Ni		NR	NR

## Hi res Extractables

		GS01SD		GS02SD		GS03SD	$\overline{}$	GS04SD		GS05SD		GS06SD		GS07SD		GS08SD	
2-Methylnaphthalene	UG/KG	0.6	J	0.9	J	33		2	J	0.8	J	_0.4	J	0.4	J	0.8 U	
Acenaphthene	UG/KG	0.5	J	1	J	100		- 5		0.9	J	0.8	Ū	0.9	U	0.8 U	_
Acenaphthylene	UG/KG	3		2		44		9		2	U	0.8	U	0.8 (	J	0.8 U	_
Anthracene	UG/KG	3		7		140		34		7		2	Ü	0.8	U	0.8 U	
Benzo(a)Anthracene	UG/KG	1	J	7		110		19		12		5	Ū	0.8	J_T	1 0	_
Benzo(b)Fluoranthene	UG/KG		NA		NA	Ţ	NA		NA		NA	,	NA	1	NA	N.	Ā
Benzo(ghi)Perylene	UG/KG		NA		NA		NA		NA		NA		NA	1	NA	N.	Ā
Benzo(k)Fluoranthene	UG/KG		NA		NA		NA		NA		NA		NA	1	NA	N.	Ā.
Benzo-a-Pyrene	UG/KG	2		4		120		17		16		4	Ü	0.8	I	0.8 U	_
Chrysene	UGAKG	1	J	10		. 200		36		26		7	Ū	1 1	J	2 U	
Dibenzo(a,h)Anthracene_	UG/KG	1	U	1	J	40		6		4		1		0.8 L	7	0.8 U	
Fluoranthene	UG/KG	2		21		190		50		34		12	U	2 (	J	2 U	_
Fluorene	UG/KG	0.5	J	2		82		6		1	J	0.3	J	0.8 (	7	0.8 U	_
Indeno (1,2,3-cd) Pyrene	UG/KG		NA		NA		NA		NA	1	NA		NA	N	VA.	N.	Ā
Naphthalene	UG/KG_	0.6	J	0.6	J	51		3	J	0.9	J	0.5	J	0.4 J		0.4 J	_
Pentachiorophenol	UG/KG	24	U	22	U	100	Ų	49	U	23	Ū	13	Ū	13 (	7	13 U	
Phenanthrene	UG/KG	2	J	12		72		11		23	Ū.	4		0.9 L	] [	0 U	
Pyrene	UG/KG	2		15	-	140		33		27		10	Ũ	2 (	7	2 U	_

Volatiles	Scan
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· · · · · · · · · · · · · · · · · · ·	i	GS01SD		GS02SD		GS03SD		GS04SD		GS05SD		GS06SD	GS07SD	GS08SD
% Molsture	%;	55		52		79		78		54		20	18	19
1,1,1-Trichioroethane	3/KG	37		34		75		70		29		13 U	12 U	14 U
	3/KQ	37		34		75		70		29		13 U	12 U	14 U
1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	3/KG	37		34	Ū	75		70		29		13 U	12 U	14 U
1,1,2-Trichloroethane	3/KG	37		34		75		70		29		13 U	12 U	14 U
1,1-Dichloroethane	3/KG	37	U	34	U	75		70		29		13 U	12 U	14 U_
1,1-Dichloroethene (1,1-Dichloroethylene)	3/KG	37	U	34	U	75	υ	70	U	29	U	13 U	12 U	14 U
1,2,3-Trichlorobenzene	3/KG		NA		NA		NA	T	NA		ÑA	NA	NA	NA NA
1,2,4-Trichiorobenzene	3/KG	37	U	34	Ū	. 75	Ū	70		29		13 U	12 U	14 U
1,2-Dibromo-3-Chioropropane (DBCP)	3/KG	37	ŪJ	34	ÜĴ	75		70		29		13 U	12 U	- 14 U
1,2-Dibromoethane (EDB)	3/KG	37	Ū	34		75		70	U	29		13 U	12 U	14 U
1,2-Dichlorobenzene	3/KG	37	U	34	Ü	75	Ū	70	Ū	29		13 U	12 U	14 U
1,2-Dichloroethane	3/KG	37	Ų	34		75		70		29		13 U	12 U	14 U
1,2-Dichioropropane	3/KG	37	Ū	34	Ū	75		70	U	29		13 U	12 U	14 U
1,3-Dichlorobenzene	3/KG	37	Ū	34	Ü	75		. 70		29		13 U	12 U	14 U
1,4-Dichlorobenzene	3/KG	37	Ū	34		75		70		29		13 U	12 U	14 U
Acetone	3/KG	94		130		350	J	520	J	51		22 J	49 J	29 J
Benzene U	3/KG	37	Ú	34	Ú.	75		70	U	29 (		13 U	12 U	14 U_
	3/KG		ÑΑ		NA		ΝĀ		NA_		NA	NA.	. NA	NA NA
	3/KG	37		34		75		70		29 (		13 U	12 U	14 U
	3/KG	37		34		75		70		29		13 U	12 U	14 U
	3/KG	37		34		75		70		29		13 U	12 UJ	14 U
	3/KG			34		51		70		29 (		13 U	12 U	14 Ú
	3/KG	37		34		75		70		29 (		13 U	12 U	14 U
	3/KG	37		. 34		75		70		29		13 U	. 12 U	14 U
	√KG			34		75		70		29 (		13 U	12 U	14 U
	KG.	37		34		75		70		29		13 U	12 U	14 U
	3/KG	37		34		75		70		29 (		13 U	12 U	14 U
	ЖG	37		34		75		70		29 1		13 U	12 U	14 U
	3/KG	37		34		75		70		29 (		13 U	12 U	14 U
	KG.	37		34		75		70		29 l		13 U	12 U	14 U
	K)	37		34		75		70		29 (		13 U	12 U	14 U
	3KG	37		34		75		70		29 t		13 U	12 U	14 U
	жg	37		34		75		70		29 (		13 U	12 U	14 U
	KG.	37		34		75		70		29 (		13 U	12 U	14 U
	XG.	37		34		75 75		70		29 t		13 U	12 U	14 U
	KG.	37		34				. 70				13 UJ	12 U	14 UJ
	VKG	37		34		75		88 70		29 t		13 UJ	12 U	14 UJ
	KG	37		34		75						13 UJ	12 U	14 UJ
	KG	37		34		75 75		70		29 (		13 U	12 U	14 U
	VKG	37 37		34 34		75		70 70		29 U		13 U 13 U	12 U 12 U	14 U
	KG	37		34		75		70		29 (			12 U	14 U
	KG KG	37		34		75		70		29 (		13 U 13 U	12 U	14 U
	/KG	37		34		75		70		29 1		13 U	12 U	14 U
	VKG	37		34		75		70		29 1		13 U	12 U	14 U
	KG	37		34		75		70		29 1				14 U
		37		34		75		70		29 (		13 U 13 U	12 U	
	VKG	37		34		75 1		70		29 (			12 U	14 U
	VKG	37		34		75 (		70		29 (		13 U	12 U	14 U
												13 U	12 U	14 U
Vinyl Chloride uc	ΚG	37	<u>-</u>	34	<u> </u>	75		70	<u> </u>	29 (	,	13 U	12 U	14 U

	!	GS01SD		GS02SD	GS03SD	GS04SD		GS05SD	GS06SD	GS07SD	GS08SD
	3/KG		NR	NR	NR	71	NJ	NR	NR	NR	NR
UNKNOWN COMPOUND	/KG		NR	NR	NR	180	J	33 J	NR	NR	8 J

#### **Data Qualiflers**

U-Analyte not detected at or above reporting limit. The number is the minimum quantitation limit. J-identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-Identification of analyte is acceptable; reported value may be blased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be blased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Table 9. Groundwater Analytical Results. Page 1 of 4

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•	[	GS01GW	GS01PB		GS02GW		GS03GW		GS04GW		GS05GW		GS06GW	4
Aluminum	UG/L	320 U.	150	S	450	J	850	_	1700	J	370	J	74	
Antimony'	UGA	2.4 U	2.4	Ú	2.4	U	2.4	C	2.4	U	2.4	U.	2.4	
Arsenic	UG/L	3.5 U	3.5	U	3.5	U	3.5	C	3.5	Ü	3.5	J	3.5	
Barlum	UG/L	20	0.6	C	120		29		26		32		0.6	U
Beryllium	UG/L	0.11	0.1	U	0.14		0.11		0.1		0.9		0.1	
Cadmium	UGAL	0.2 U	0.2	C	0.2		0.2	U	0.2	U	0.2	υ	0.2	
Calcium	UG/L	4600	430		16000		3600		3500		15000		430	
Chromium	UG/L	1.8 R	0.6		0.92		1.8		1.6		0.6		0.6	
Cobalt	UG/L	1.2	0.9	ט	1.2		0.9		0.9		10		0.9	
Copper	UG/L	1.5 U	0.6	ح	0.73		4.1		0.78		0.6		0.67	
Cyanide	UGAL	N/	Α	NA		NA		NA		NA	1	NA		NA
Iron	UG/L	1500	52		1400		2100		1900		12000		52	
Lead	UG/L	2 U		כ		U		U		U		Ü		U
Magnesium	UG/L	1500	15		5500		940		930		8200		15	
Manganese	UG/L	23	0.5		160		110		30		320		0.5	
Nickel	UG/L	1.5 U	1.5		3.5		1.5		3.3		20		1.5	
Potassium	UG/L	520	130		1500		870		1800		1900		100	
Selenium	UG/L	2.4 U	2.4		2.4		2.4		2,4		2.4		2.4	
Silver	UG/L	0.7 U	0.7		0.7		0.7	U	0.7		0.7		0.7	
Sodium	UG/L	40000	410		120000		29000		5000		86000		400	
Thallium	UG/L	3.1 U	3.1		3.1		3.1		3.1		3.1		3.1	
Total Mercury	ŲGΛL	0.1 U	0.1		0.1		0.1	U	0.1		0.1		0.1	
Vanadium	ÜGΛL	0.6 U	0.6		0.82		2		1.9		0.6	U	0.6	
Zinc	ŲG/L	11 U	11	U	13		18		11	U	. 42		11	U

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		GSUIGW	GSOIPB	GS02GW	G503GW	GS04GW	GSUSGW	GSOGGW
4,4'-DDD (p,p'-DDD)	ψG/L	0.02 UJ	NR	0.02 UJ	0.02 UJ	0.02 UJ	0.02 UJ	NR
4,4'-DDE (p,p'-DDE)	ÜG/L	0.02 UJ	NR	0.02 UJ	0.02 UJ	0.02 UJ	0.02 UJ	NR
4,4'-DDT (p,p'-DDT)	UGA	0.02 UJ	NR	0.02 UJ	0.02 UJ	0.02 UJ	0.02 UJ	NR
Aldrin	UG/L	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
alpha-BHC	UG/L	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
alpha-Chiordane /2	UG/L	0.01 U	NA	0.01 U	0.01 U	0.01 U	0.01 U	NR
beta-BHC	UGAL	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
delta-BHC	UG/L	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
Dieldrin	UG/L	0.02 U	NR	0.02 U	0.02 U	0.02 U	0.02 U	NR
Endosulfan I (alpha)	UGAL	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
Endosulfan İl (beta)	UG/L	0.02 U	NR NR	0.02 U	0.02 U	0.02 U	0.02 U	NR
Endosulfan Sulfate	UG/L	0.02 U	NR	0.02 U	0.02 U	0.02 U	0.02 U	NR
Endrin	ng/r	0.02 UJ	NR	0.02 UJ	0.02 UJ	0.02 UJ	0.02 UJ	NR
Endrin Aldehyde	UG/L	0.02 UJ	NR	0.02 UJ	0.02 UJ	0.02 UJ	0.02 UJ	NR
Endrin Ketone	UG/L	0.02 UJ	NR	0.02 UJ	0.02 UJ	0.02 UJ	0.02 UJ	NR
gamma-BHC (Lindane)	UG/L	. 0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR.
gamma-Chlordane /2	UGAL	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
Heptachlor	UG/L	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
Heptachlor Epoxide	UG/L	0.01 U	NR	0.01 U	0.01 U	0.01 U	0.01 U	NR
Methoxychlor	lJG/L	0.1 U	NR	0.1 U	0.1 U	0.1 U	0.1 U	NR
PCB-1016 (Aroclor 1016)	UG/L	0.2 U	NR .	0.2 U	0.2 U	0.2 U	0.2 U	NR
PCB-1221 (Aroclor 1221)	UGAL	0.4 U	NR	0.4 U	0.4 U	0.4 U	0.4 U	NR
PCB-1232 (Aroclor 1232)	UG/L	0.2 U	NR	0.2 U	0.2 U	0.2 U	0.2 U	NR
PCB-1242 (Aroclor 1242)	UG/L	0.2 U	NR	0.2 U	0.2 U	0.2 U	0.2 U	NR
PCB-1248 (Arocior 1248)	UG/L	0.2 U	NR	0.2 U	0.2 U	0.2 U	0.2 U	NR
Pasticidas & Arociora Scan			<del></del>					

	<u>i</u> .	GSUIGW	GSUIPB	GS02GW	GSU3GW	GS04GW	GSUSGW	GSUGGW
PCB-1254 (Arocior 1254)	UG/L	0.2 U	NR	0.2 U	0.2 U	0.2 U	0.2 U	NR
PCB-1260 (Arocior 1260)	UG/L	0.2 U	NR	0.2 U	0.2 U	0.2 U	0.2 U	NR
Toxaphene	UG/L	1 0	NR	1 U	1 U	1 U	1 U	NR

## Extractables Scan

		GS01GW	GS01PB	GS02GW	GS03GW	GS04GW	GS05GW	GS06GW
(3-and/or 4-)Methylphenol	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
1,1-Biphenyl	UG/L	5 UJ	NR	5 UJ	5 UJ	5 U	5 U	NP
1,2,4,5-Tetrachiorobenzene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NA
2,4,5-Trichlorophenoi	UG/L	20 UJ	NR	20 UJ	20 U	20 U	20 U	NR
2,4,6-Trichlorophenol	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NA
2,4-Dichlorophenol	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NA
2,4-Dimethylphenol	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NA
2,4-Dinitrophenol	UGAL	20 UJ	NR	20 UJ	20 UJ	20 UJ	20 UJ	NA
2,4-Dinitrotoluene	UG/L_	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
2,6-Dinitrotoluene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NF
2-Chloronaphthalene	UG/L	5 UJ	NR NR	5 UJ	5 U	5 U	5 U	NA NA
2-Chlorophenol	UG/L UG/L	20 UJ	NR NR	20 UJ	5 U 20 UJ	5 U 20 UJ	20 UJ	NF
2-Methyl-4,6-Dinitrophenol 2-Methylnaphthalene	UG/L	5 UJ	NR NR	5 UJ	5 U	20 U3	5 U	NF
2-Methylphenol	UG/L	5 UJ	NR	5 UJ	5 U	. 5 U	5 U	NF
2-Nitroaniline	UG/L	20 UJ	NR	20 UJ	20 U	20 U	20 U	NF
2-Nitrophenol	UG/L	5 UJ	NR NR	5 UJ	5 0	5 U	5 U	NF
3,3'-Dichlorobenzidine	UG/L	5 UJ	NR	5 UJ	5 UJ	5 UJ	5 0	NR
3-Nitroaniline	UG/L	20 UJ	NR	20 UJ	20 U	20 U	20 U	NR
4-Bromophenyl Phenyl Ether	UGAL	5 UJ	NR	5 UJ	5 U	5 U	5 U	NA
4-Chloro-3-Methylphenol	UG/L	5 UJ	NR	5 03	5 0	5 U	5 U	NA
4-Chioroaniline	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
4-Chlorophenyl Phenyl Ether	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
4-Nitroanlline	UG/L	20 UJ	NR	20 UJ	20 U	20 U	20 U	NR
4-Nitrophenoi	UG/L	20 UJ	NR	20 UJ	20 U	20 U	20 U	NR
Acenaphthene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Acenaphthylene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NA
Acetophenone	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Anthracene	UG/L	5 UJ 5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Atrazine	UG/L	5 UJ	NR NR	5 UJ	5 U	5 U	5 U	NR
Benzaldehyde Benzo(a)Anthracene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR NR
Benzo(b)Fluoranthene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR NR
Benzo(ghi)Perylene	UGAL	5 UJ	NR	5 UJ	5 0	5 U	5 U	NR.
Benzo(k)Fluoranthene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Benzo-a-Pyrene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Benzyl Butyl Phthalate	UGAL	5 UJ	. NR	5 UJ	5 UJ	5 U	5 Ü	NR
bis(2-Chloroethoxy)Methane	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
bis(2-Chloroethyl) Ether	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
bis(2-Chloroisopropyl) Ether	UG/L	5 UJ	NR	5 UJ	5 U	5 U	. 5 U	NR
bis(2-Ethylhexyl) Phthalate	UG/L	5 UJ	NR	5 UJ	5 UJ	5 U	5 U	NR
Caprolactam	UG/L	5 UJ	NR	5 UJ	5 UJ	5 U	5 U	NR
Carbazole	UG/L	NA NA	NR	NA NA	NA	NA NA	NA NA	NR
Chrysene	UG/L UG/L	5 UJ	NR NR	5 UJ	5 U	5 U	5 U	NR
Dibenzo(a,h)Anthracene Dibenzofuran	UG/L	5 UJ	NR NR	5 UJ 5 UJ	5 U	5 U	5 U 5 U	NR NR
Diethyl Phthalate	UG/L	5 UJ	NR	5 UJ 5 UJ	5 UJ	5 U	5 U	NR NR
Dimethyl Phthalate	UG/L	5 UJ	NR	5 UJ	5 UJ	5 U	5 0	NR NR
Di-n-Butyiphthalate	UGAL	5 UJ	NR	5 UJ	5 UJ	5 0	5 0	NR.
Di-n-Octylphthalate	UG/L	5 UJ	NR	5 UJ	5 UJ	5 0	5 U	NR.
Fluoranthene	UGAL	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Fluorene	UG/L	5 UJ	. NR	5 UJ	5 U	5 0	5 U	NR
Hexachlorobenzene (HCB)	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Hexachlorobutadiene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Hexachlorocyclopentadiene (HCCP)	UG/L	5 UJ	NR	5 UJ	5 UJ	5 ŪJ	5 UJ	NR
Hexachloroethane	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR.
Indeno (1,2,3-cd) Pyrene	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR
Isophorone	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	. NR
Naphthalene	UG/L	5 UJ	NR NR	5 UJ	5 U	5 U	5 U	NR
Nitrobenzene	UGAL	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR.
n-Nitrosodi-n-Propylamine	UG/L	5 UJ	NR	5 UJ	5 U	5 U	5 U	NR

Extractables Scan								
	<u> </u>	GS01GW	GS01PB	GS02GW	GS03GW	GS04GW	GS05GW	GS06GW
n-Nitrosodiphenylamine/Diphenylamine	ŲG/L	5 UJ	NR	5 UJ		5 U	5 U	N
Pentachlorophenol	ÜGΛ	5 UJ	NR	5 ÚJ		5 U	5 U_	N
Phenanthrene	UG/L	5 UJ	NR	5 UJ		5 U	5 U	N
Phenol	UG/L	5 UJ	NR	5 UJ		5 U	5 U	N
Pyrene	UGAL	5 UJ	NR	5 UJ	5 U	5 U	5 U	N
Miscellaneous Extractables		GS01GW	<del></del>	GS02GW	GS03GW	<del></del>	GS05GW	<del> </del>
2 UNKNOWN COMPOUNDS	UGA.	NR		GSUZGW NF			4.7 J	<del> </del>
D-LIMONENE	UGAL	2,2 NJ	·	NF		<del></del>	NF.	
LIMONENE	UG/L	NR		2.2 NJ			NF	
Volatiles Scan		,						Leanny
# 4 4 Talahla mathana	1104	GS01GW	GS01PB	GS02GW	GS03GW 0.5 U	GS04GW	GS05GW	GS06GW
1,1,2,2-Tetrachioroethane	UG/L UG/L	0.5 U	NR NR	0.5 U	0.5 U	0.5 U	0.5 U	NI NI
1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	UGAL	0.5 U	NR.	0.5 U	0.5 U	0.5 U	0.5 U	· Ni
1,1,2-Trichloroethane	LGA	0.5 U	NR NR	0.5 U	0.5 U	0.5 U	0.5 U	N.
1,1-Dichloroethane	UGAL	0.5 U	NR NR	0.5 U	0.5 U	0.5 U	0.5 U	N
1,1-Dichloroethene (1,1-Dichloroethylene)	UG/L	0.5 U	NR NR	0.5 U	0.5 U	0.5 U	0.5 U	NI NI
1,2,3-Trichlorobenzene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
1,2,4-Trichlorobenzene	DGZ	0.5 U	NA	0.5 U	0.5 U	0.5 U	0.5 U	NI NI
1,2-Dibromo-3-Chloropropane (DBCP)	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
1,2-Dibromoethane (EDB)	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
1,2-Dichlorobenzene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
1,2-Dichloroethane	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
1,2-Dichloropropane	UGAL	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
1,3-Dichlorobenzene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	· N
1,4-Dichlorobenzene	UG/L	0.5 U	NR.	0.5 U	0.5 U	0.5 U	0.5 U	N
Acetone	UG/L	5 U	NR	5 U	5 U	5 U	5 U	N
Benzene	UGAL	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Bromochloromethane	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Bromodichioromethane	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
Bromoform	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Bromomethane	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
Carbon Disulfide	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Carbon Tetrachioride	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Chlorobenzene	UG/L	0.5 U	NA	0.5 U	0.5 U	0.5 U	0.5 U	N
Chloroethane	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NF
Chloroform	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NF
Chloromethane	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
cis-1,2-Dichloroethene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
cis-1,3-Dichloropropene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
Cyclohexane	UG/L	0.5 U	NR NR	0.5 U	0.5 U	0.5 U	0.5 U	NF
Dibromochioromethane	UGAL	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NF
Dichlorodiffuoromethane	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NF
Ethyl Benzene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NF
Isopropylbenzene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NF NF
Methyl Acetate	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NF
Methyl Butyl Ketone	UG/L	5 U	NR NR	5 U	5 U	5 U	5 U	NF NF
Methyl Ethyl Ketone	UG/L	6.1 UJ	NR	5 01	6.9 UJ	7.2 J	6.6 UJ	
Methyl Isobutyl Ketone Methyl T-Butyl Ether (MTBE)	UG/L UG/L	5 U	NR NR	5 U 0.5 U	5 U	5 U 0.5 U	5 0	NF NF
Methyl 1-Butyl Ether (MTBE) Methylcyclohexane	UG/L	0.5 U	NR NR	0.5 U	0.5 U	0.5 U	0.5 U	Ni Ni
Methylene Chloride	UG/L	0.5 U	NR NR		0.5 U	0.5 U		N
Styrena	UG/L	0.5 U	NR NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Tetrachloroethene (Tetrachloroethylene)								
·	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Foluene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	N
Total Xylenes	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NI
rans-1,2-Dichloroethene	: UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	Ni
rans-1,3-Dichloropropene	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	l Ni

#### Volatiles Scan

<u> </u>		GS01GW	GS01PB	GS02GW	GS03GW	GS04GW	GS05GW	GS06GW
Trichloroethene (Trichloroethylene)	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NR
Trichlorofluoromethane (Freon 11)	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NA
Vinyl Chioride	UG/L	0.5 U	NR	0.5 U	0.5 U	0.5 U	0.5 U	NR

#### Miscellaneous Volatiles

	:	GS01GW	GS02GW	GS03GW	/	GS050	3W
INDANE	UG/L	NR	N	R 1.3	NJ		NR

#### Data Qualifiers

U-Analyte not detected at or above reporting limit. The number is the minimum quantitation limit.

J-Identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-identification of analyte is acceptable; reported value may be blased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be blased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolities of technical chlordane

Table 10. Surface Soil Miscellaneous Analytical Results. Page 1 of 8

MISCEITATIOUS EXTRACTABLES	Units	GS01SS		GS02SS	GS03SS	GS04SS		GS05SS		GS06SS	GS07SS
[3,4:9,10] DIBENZPYRENE	UG/KG		NR	NR NR	NR NR	430433	NR	180	N.I	NR NR	NR
1-DOCOSENE	UG/KG		NR	NR	NR.	<del> </del>	NR	180	NR	NR.	NR NR
1,2:3,4-DIBENZPYRENE	UG/KG		NR	NR	NR NR	<del>  </del>	NR	<del> </del>	NR	NR NR	NR NR
1,2:4,5-DIBENZPYRENE	UG/KG		NR	NR.	NR NR	<del> </del>	NR	<del> </del> -	NR	NR NR	NR
1,2:7,8-DIBENZPHENANTHRENE	UG/KG		NR	NR NR	NR NR	+	NR	260	NJ	NR NR	NR NR
1,2:7,8-DIBENZPHENANTHRENE	UG/KG		NA	NR NR	NR.	<del>- </del> -	NR	200	NR	NR NR	NR
10 UNKNOWN COMPOUNDS	UG/KG		NR	NR NR	NR NR		NR	<del></del>	NR	NR NR	NR
10 UNKNOWN PAHS	UG/KG		NR	NR NR	NR NR	<del> </del> -	NR	<del> </del>	NR	NR NR	NR NR
11 UNKNOWN COMPOUNDS	UG/KG		NR	NR	NR.	<del> </del>	NR	2700	J	NR NR	NR
11H-BENZO [A] FLUORENE	UG/KG		NR	NR NR	NR.		NR			NR NR	NR NR
11H-BENZO [B] FLUORENE	UG/KG		NR	NR NR	- NR		NR	4/0	NR	NR NR	740 NJ
12 UNKNOWN COMPOUNDS	UG/KG	1400	INIT	NR NR	NR.	<del></del>	NR	<del></del>	NR	NR NR	NR NR
14 UNKNOWN COMPOUNDS	UG/KG		NR	NR.	NR NR	<del>                                     </del>	NR	<del> </del>	NR	2800 J	NR.
15 UNKNOWN COMPOUNDS	UG/KG		NR	NR NR	NR NR	<del></del>	NR	<del> </del>	NR	2800 J NR	NR NR
16 UNKNOWN COMPOUNDS	UG/KG		NR	NR	NR NR	<del></del>	NR NR		NR	NR NR	NR NR
17 UNKNOWN COMPOUNDS	UG/KG		NR	NR NR	NR NR	<del> </del>	NR NR	<del> </del>	NR	NR NR	NR NR
	UG/KG		NA	NR NR				<del> </del>		NA NA	NR NR
19 UNKNOWN COMPOUNDS 1-EICOSANOL	UG/KG		NR NR	NR NR	NR NR		NR NR	<del> </del>	NR NR	NR NR	NR NR
1-HEPTADECENE	UG/KG		NR	NR NR	NR NR		NH NR	+	NR	120 NJ	NR NR
1-OCTADECENE	UG/KG		NR NR	NR NR	NR NR		NR NR		NR NR	120 NR	NR NR
1-TETRADECENE	UG/KG		NR	NR NR	NR NR		NR		NR	NR NR	NR NR
2 UNKNOWN KETONES	UG/KG		NR I	NR NR	NR.	250			NR	NR NR	NR NR
2.2'-BINAPHTHALANE	UG/KG		NR	NR NR	NR NR		NR		NR.	NR NR	NR NR
2.2'-BINAPHTHALENE	UG/KG		NR	NR	NR.		NR		NR	NR NR	NR.
21 UNKNOWN COMPOUNDS	UG/KG		NR	NR NR	NR.		NR		NR	NR NR	NR NR
25 UNKNOWN COMPOUNDS	UG/KG		NR	. NR	NR NR	+	NR		NR	NR NR	NR NR
3 UNKNOWN COMPOUNDS	UG/KG		NR	NR.	NR.		NR		NR	NR NR	NR NR
3 UNKNOWN PAHS	UG/KG		NR	NB	NR.		NR		NB	NR I	NR NR
3,4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG		NR	NR	NR		NR		NR	NR NR	NR.
4 UNKNOWN COMPOUNDS	UG/KG		NR	NR	NR		NR		NR	NR NR	NR
4 UNKNOWN PAHS	UG/KG		NR	NR	NR	1	NR	. 1500	.1	NR	NR
5 UNKNOWN COMPOUNDS	UG/KG		NR	2200 J	NR		NR		NR	NR NR	NR
5, 12-NAPHTHACENEDIONE	UG/KG		NR	NR	NR		NR		NR	NR	NR NR
5,12-NAPHTAHCENEDIONE	UG/KG		NR	. NR	NR		NA		NA	NR NR	NR
5.12-NAPHTHACENEDIONE	UG/KG		NR	NR NR	NR.		NR	330		NR NR	NR NR
6 UNKNOWN COMPOUNDS	UG/KG		NR	NR	NR		NR-		NR	NR.	NR
6 UNKNOWN PAHS	UG/KG		NR	NR	NR		NR		NR	NR	NR NR
7 UNKNOWN COMPOUNDS	UG/KG		NR	NR	NR	2000	1		NR	NR NR	NR NR
7 UNKNOWN PAHS	UG/KG		NR	NR	NR		NR		NR	NR I	5300 J
7H-BENZ [DE] ANTHRACEN-7-ONE	UG/KG		NR	NR	NR		NR	310		NR	850 NJ
8 UNKNOWN COMPOUNDS	UG/KG		NR	NR	NR		NR		NR	NR NR	5900 J
8 UNKNOWN PAHS	UG/KG		NR	NR	NR		NA		NR	NR	NR
9, 10-ANTHRACENEDIONE	UG/KG		NR	NR	NR		NR.		NJ	NR	NR
9,10-ANTHRACENEDIONE	UG/KG		NR	NR	NR		NR		NR	NR	890 NJ
9H-FLUOREN-9-ONE	UG/KG		NR	NR	NR		VR.		NR	NR	460 NJ
9-OCTADECENE, (E) -	UG/KG		NA	NR	NR		VR.		NR	170 NJ	NB NB
ANTHRACENE, 2-METHYL-	UG/KG		NR.	NR-	NR		VR.		NR	NR	1000 J
BENZ (A) ANTHRACENE-7, 12-DIONE	UG/KG	1	NR	NR	NR		VR.		NR	NR	NR
BENZ (A) ANTHRACENE-7,12-DIONE	UG/KG	-	VR.	NR	NR		VR.		NR	NR	540 NJ
BENZ [A] ANTHRACENE, 7-METHYL-	UG/KG		VR	NR	NR		VR.		NR	NR	NR
BENZ [E] ACEPHENANTHRYLENE	UG/KG	,	NA	NR	NR		NR.		NR	NR	NR
BENZO [B] NAPHTHO [2,1-D] THIOPH	UG/KG		VR	NR	NR		VA.		NR	NR	NR
BENZO [B] NAPHTHO [2,1-D] THIOPHENE	UG/KG	1	VR.	NR	NR		VR.		NJ	NA	NR
BENZO [B] NAPHTHO [2,1-D] THIPHENE	UG/KG		NR	NR	NR		NR.		NR	NR	670 NJ
BENZO [B] NAPHTHO [2,3-D] FURAN	UG/KG		VR	NR	NR		VR.		NR	NR	NB NB
BENZO [B] TRIPHENYLENE	UG/KG		VR.	NR	NR		VR.		NR	NR	NR:
BENZO [E] PYRENE	UG/KG		VR.	NR	NR	<del></del>	VR.	290		NR NR	1200 NJ
		<del></del>	لمنتب		1,,(1	<u></u>	** *	200		1,41,1	1200 110

	Units	GS01SS		GS02SS		GS03SS	_	GS04\$\$		GS05SS		GS06SS		GS07SS	$\overline{}$
BENZO [J] FLUORANTHENE	UG/KG	!	NR		NR	1	NR		NR		NR		NR	N	VR
BENZOFLUORENE	UG/KG		NR		NR		NR		NR	160	J		NR	I N	VR.
CARYOPHYLLENE	UG/KG		NR		NR		NR		NR		NR		NR		VR.
CHOLESTEROL	UG/KG		NR		NR	1	NR		NR		NR	T	NR	N	VR
CHRYSENE, 1-METHYL-	UG/KG		NR		NR	1	NR		NR		NR		NR	N	VR.
CHRYSENE, 5-MEHTHYL-	UG/KG		NR .		NR	1	NR		NR		NR		NR	580 N	۱IJ
CHRYSENE, 5-METHYL-	UG/KG	1.	NR		NR	1	NR		NA	1	NR		NR	10	VR.
CHRYSENE, 6-METHYL-	UG/KG		NR		NR		NR		NR	270	Ŋ		NR	N	VR.
CYCLOPENTA (DEF) PHENANTHRENONE	UG/KG		NR		NR	1	NR	]	NR		NR		NR	N	NR
D:C-FRIEDOOLEANAN-3-ONE	UG/KG		NR		NR	1	ŇR		NR		NR	860	NJ	N	VR.
HEXADECANOIC ACID	UG/KG		NR		NR	1	NR		NR		NR	100	NJ	N	NR.
METHYLANTHRACENE	UG/KG	-:	NR		NR	,	NR		NR	190	J		NR	N	VR.
NAPHTHALENE, 2-PHENYL-	UG/KG		NR		NR		NR		NR		NR		NR .	760 N	1 <u>J</u>
PENTADECANOIC ACID	UG/KG		NR		NR	1	NR		NR		NR		NR	N	VR.
PERYLENE	UG/KG		NR		NR	1	NR		NR		NR		NR	N	NR.
PHENANTHRENE, 1-METHYL-	UG/KG		NR		NR	I	NR		NR		NR		NR	١	VR.
PHENANTHRENE, 2,3-DIMETHYL-	UG/KG	7	NR		NR	1	NR		NR	160	NJ:		NR	I N	NR.
PHENANTHRENE, 2-METHYL-	UG/KG	1	NA		NR	1	NR		NR		NR		NR	870 N	IJ.
PHENANTHRENE, 4,5-DIMETHYL-	UG/KG		NR		NR		NR		NR		NR		NR	830 N	۱ <u>۱</u>
PYRENE 2-METHYL-	UG/KG		NR		NR	1	NR		NR		NR		NR	1200 N	IJ
PYRENE, 1,3-DIMETHYL-	UG/KG	-	NR		NR	1	NR		NR	280	NJ		NR	N	iR
PYRENE, 1-METHYL-	UG/KG		NR		NR	1	NR		NR		NR		NR	N	NR.
PYRENE, 2-METHYL-	UG/KG		NA		NR	1	NR		NR	250	NJ		NR	N	VR.
PYRENE, 4-METHYL-	UG/KG	1	NR		NR	1	NR.		NR		NR		NR	740 N	1J
STIGMAST-4-EN-3-ONE	UG/KG		NR	190	NJ	1	NR	350	J		NR .	460	NJ	N	IR.
TESTOSTERONE	UG/KG		NR		NR	1	NR		NR		NR		NR		VR_
TRIPHENYLENE, 2-METHYL-	UG/KG		NR		NR		NR		NR		NR		NR	N	1R
UNKNOWN ALCOHOL	UG/KG		NR		NR		NR	92	J		NR		NR		IR .
UNKNOWN ALKENE	UG/KG	1	NR		NR	1	NR		NR		NR		NR	N	NR.
UNKNOWN CARBOXYLIC ACID	UG/KG		NR		NR		VR.		NR:	<u> </u>	NR		NR		iR
UNKNOWN COMPOUND	UG/KG		NR		NR	91 J			NR		NR		NR		iR
UNKNOWN PAH	UG/KG		NR		NR		NR_		NR		NR		NR		IR
VALENCENE	UG/KG		NR		NR	l N	VR.		NR		NR		NR	N	IR.

		GS01SS		G\$02SS		GS03SS		GS04SS		GS05SS	GS06SS		GS07SS
2 UNKNOWN COMPOUNDS	UG/KG	19	J	51	J	N	IR .		NR	NR	20	5	NR
3 UNKNOWN COMPOUNDS	UG/KG	-	NR		NR	N	IR	-	NR	NR		NR	NR
ACETALDEHYDE	UG/KG	1	NR		NR	N	IR :		NR	NR		NR	NR
BUTANAL	UG/KG		NR		NR	N	IR		NR	NR		NR	NR
HEXANAL	UG/KG		NR		NR	N	IR		NR	NR	64	NJ	NR
PENTANAL	UG/KG		NR		NR	N	IR		NR	NR	T	NR	NR
PROPANAL, 2-METHYL-	UG/KG	!	NR		NR	. N	IR		NR	NR		NR	NR
UNKNOWN COMPOUND	UG/KG		NR		NR	N	R	19	J	7 J		NR	. 10 J
UNKNOWN COMPOUNDS	UG/KG		NR		NR	6 J			NR	NR	· · · · ·	NR	NR

Table 10. Surface Soil Miscellaneous Analytical Results. Page 3 of 8

	Units	GS08SS		GS0955		GS10SS	T-	GS11SS	GS12SS	GS13SS		GS14SS	
[3,4:9,10] DIBENZPYRENE	UG/KG	1	NA		NR		NR	NI	NR NR		NR		NR
1-DOCOSENE	UG/KG		NR		NR	1	NR	N	NR NR		NR		NR
1,2:3,4-DIBENZPYRENE	UG/KG		NR		NR		NR	NE	NR NR		NR		NR
1,2:4,5-DIBENZPYRENE	UG/KG	880	NJ		NR		NR	NI	NR NR	-	NR		NR
1,2:7,8-DIBENZPHENANTHRENE	UG/KG	:	NR		NR		NR	N	NR NR		NA	550	NJ
1.2:7.8-DIBENZPHENANTHRRENE	UG/KG		NR		NR		NR	480 J	NR		NR		NR
10 UNKNOWN COMPOUNDS	UG/KG		NR	1900	J		NR	Ni	NR NR		NR		NR
10 UNKNOWN PAHS	UG/KG	:	NR		NR	9400	J	NE	NR.	1	NR		NR
11 UNKNOWN COMPOUNDS	UG/KG		NA		NR	1	NR	N	NR NR		NR		NR
11H-BENZO [A] FLUORENE	UG/KG		NR		NR	2200	NJ	NI	NR NR		NA	1700	NJ
11H-BENZO (B) FLUORENE	UG/KG		NR	· · · · · · · · · · · · · · · · · · ·	NR		NR	N/	NR NR	1100	NJ		NR
12 UNKNOWN COMPOUNDS	UG/KG	-	NR		NR		NR	3600 J	NR	19000	J		NR
14 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR	N	NR.		NR		NR
15 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR	N	NR.		NR	1 1	NR
16 UNKNOWN COMPOUNDS	UG/KG	<b> </b>	NR		NR	<del> </del>	NR	NI		<del> </del>	NR		NR
17 UNKNOWN COMPOUNDS	UG/KG	23000	J		NR	1	NR	N		+	NR		NR
19 UNKNOWN COMPOUNDS	UG/KG		NB		NR		NR	N		1	NR		NR
1-EICOSANOL	UG/KG		NR		NR		NR	N			NR		NR
1-HEPTADECENE	UG/KG	1	NR	230	ΝJ		NR	NI		<del>                                     </del>	NR	· · · · ·	NR
1-OCTADECENE	UG/KG	1	NR	230			NR	N			NR		NR
1-TETRADECENE	UG/KG	1	NR	140	NJ	<del> </del>	NR	N	NR NR	+	NR	1	NR
2 UNKNOWN KETONES	UG/KG	1	NR		NR	1	NR	NF		1	NR		NR
2,2'-BINAPHTHALANE	UG/KG	1	NR.		NR		NR	N N	NR NR	560	NJ		NR
2,2'-BINAPHTHALENE	UG/KG		NR		NR	560	ÑJ	NF	NR NR	1	NR		NR
21 UNKNOWN COMPOUNDS	UG/KG	·	NR_		NR	1	NR	NF	NR.	1.	NR	. 1	NR
25 UNKNOWN COMPOUNDS	UG/KG	1	NR		NR		NR	NF	NR.		NR		NR
3 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR	NF	410 J		NR		NR
3 UNKNOWN PAHS	UG/KG		NR		NR		NR	NF			NR_		NR
3,4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG		NR		NR			NF			NR		NJ
4 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR	NF			NR		NR
4 UNKNOWN PAHS	UG/KG	5400			NR		NR	NF			NA		NR
5 UNKNOWN COMPOUNDS	UG/KG		NR_		NR		NR	NF			NR		NR
5, 12-NAPHTHACENEDIONE	UG/KG		NR		NR		NR	NF			NR		NR
5,12-NAPHTAHCENEDIONE	UG/KG		NR		NR		NR	NF			ŊĴ		NR
5,12-NAPHTHACENEDIONE	UG/KG	1400			NR	690	NJ	NF.			NR	380	
6 UNKNOWN COMPOUNDS	UG/KG		NR		NR	4100	J	NF			NR		NR
6 UNKNOWN PAHS	UG/KG	<u> </u>	NR		NR		NR	2100 J	NR	3233	J		NR
7 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR	NF			NR		NR
7 UNKNOWN PAHS	UG/KG	1	NR		NR		NR	NF			NR		NR_
7H-BENZ [DE] ANTHRACEN-7-ONE	UG/KG	910			NR	1100		430 NJ	NR	1500			NJ
8 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR	NF			NR N	5500	
8 UNKNOWN PAHS	UG/KG		NR		NR		NR	NF			NR	5600	
9, 10-ANTHRACENEDIONE	UG/KG	1 7	NR		NR	660		260 NJ	NR		NR		NR
9,10-ANTHRACENEDIONE	UG/KG		NR		NR		NR .	NF	NR		Ŋ		NR
9H-FLUOREN-9-ONE	UG/KG		NR		NR		NR	NF	NR		3		NR
9-OCTADECENE, (E) -	UG/KG		NR		NR		NR	NF			NR		NR
ANTHRACENE, 2-METHYL-	UG/KG		NR		NR		NR	NF	NR		NR .		NR
BENZ (A) ANTHRACENE-7, 12-DIONE	UG/KG		NR		NR		NR	NF			NR.		NR_
BENZ (A) ANTHRACENE-7,12-DIONE	UG/KG	1 .	NR		NR		NR	330 NJ	NR		NR		NR
BENZ [A] ANTHRACENE, 7-METHYL-	UG/KG		NR		NR		NR	NF			NR ·		NR
BENZ [E] ACEPHENANTHRYLENE	UG/KG		NR		NR		NR	NF	NR		NR		NR
BENZO [B] NAPHTHO [2,1-D] THIOPH	UG/KG		NR		NR		NR	NF	NR		NR		NR
BENZO [B] NAPHTHO [2,1-D] THIOPHENE	UG/KG		NR		NR		NA	NF	NR		NR		NJ
BENZO [B] NAPHTHO [2,1-D] THIPHENE	UG/KG		NR		NR		NR	NF	NR		VR.		NR
BENZO [B] NAPHTHO [2,3-D] FURAN	UG/KG		NR		NR		NR	NF			M.		NR
BENZO [B] TRIPHENYLENE	UG/KG		NR		NR		NR	380 NJ	NR		NE.		NR
BENZO [E] PYRENE	UG/KG		NR		NR	3300	NJ	1700 NJ	NR	1 1	VR	2600	NJ

Miscontinuous Extractables	Units	GS08SS		GS09SS		GS10SS		GS11SS		GS12SS	GS13SS		GS14SS
BENZO [J] FLUORANTHENE	UG/KG	-	NR		IR		NR		NR	NR I		NR.	NR
BENZOFLUORENE	UG/KG	920	J	N	VA		NR	l ———	NR	NR		NR	. NR
CARYOPHYLLENE	UG/KG		NR	l N	JR		NR		NR	NR		NR	NR
CHOLESTEROL	UG/KG		NR	N	IR		ŇR		NR	NR		NR	NR
CHRYSENE, 1-METHYL-	UG/KG	1200	NJ	N	JR		NR		NR	NR.		NR	NR
CHRYSENE, 5-MEHTHYL-	UG/KG		NR	N	IR		NR		NR	NR ·		NR	NR
CHRYSENE, 5-METHYL-	UG/KG		NR	N	1R		NR		NR	NR	910	NJ	NR
CHRYSENE, 6-METHYL-	UG/KG	:	NR	l N	√R		NR		NR:	NR		NR	NR
CYCLOPENTA (DEF) PHENANTHRENONE	UG/KG	:	NR	N	1R	940	NJ	320	NJ	NR	1400	NJ	470 NJ
D:C-FRIEDOOLEANAN-3-ONE	UG/KG		NR	l N	1R		NR		NR	NR NR		NR	NR
HEXADECANOIC ACID	UG/KG		NR	l N	IR -		NR	670	ŊJ	NR NR		NR	NR
METHYLANTHRACENE	UG/KG		NR		1B		NR		NR	NR NR		NR	NR.
NAPHTHALENE, 2-PHENYL-	UG/KG		NR	l N	1R		NR		NR	NR		NR	NR
PENTADECANOIC ACID	UG/KG		NR	97 N	IJ		NR		NR	NR		NR	NR
PERYLENE	UG/KG	3900	NJ		1R	1000	NJ		NR	NR		NR	710 NJ
PHENANTHRENE, 1-METHYL-	UG/KG		NR	N	1R		NR		NR	NR	1100	NJ	NR
PHENANTHRENE, 2,3-DIMETHYL-	UG/KG		NR	l N	IR ]		NR		NR	NR	1300	NJ	NR
PHENANTHRENE, 2-METHYL-	UG/KG		NR	l N	IR		NR		NR	NR		NR	NR
PHENANTHRENE, 4,5-DIMETHYL-	UG/KG		NR		NR		NR		NR .	NR		NR	NR
PYRENE 2-METHYL-	UG/KG		NR	l N	IR.		NR		NR	. NR		NR	NR
PYRENE, 1,3-DIMETHYL-	UG/KG		NR	L N	IA		NR		NR	NR		NR	NR
PYRENE, 1-METHYL-	UG/KG	880			ŧ₽ Ţ	710		390		NR	1400	Z	1300 NJ
PYRENE, 2-METHYL-	UG/KG	1000		N	IR	1700	NJ		NR	NR	660	Ŋ	720 NJ
PYRENE, 4-METHYL-	UG/KG		R		IR		NR		NR	NR		NR	NR
STIGMAST-4-EN-3-ONE	UG/KG		NR	N			NR		NR	NR		NR	NR
TESTOSTERONE .	UG/KG		NR		IR [		NR		NR	NR		NR	NR
TRIPHENYLENE, 2-METHYL-	UG/KG		NR		IR	800		370		NR	1	NR	840 NJ
UNKNOWN ALCOHOL	UG/KG		NR		IR		NR		NR	NR		NR	NR
UNKNOWN ALKENE	UG/KG		NR		IR		NR_		NR .	NR		NR	NR
UNKNOWN CARBOXYLIC ACID	UG/KG		NR		IR	720	J		NR	NR		NR	NR
UNKNOWN COMPOUND	UG/KG		NR		R		NR		NR	NR		NR	NR
UNKNOWN PAH	UG/KG		NR .		R		ВZ		NR	NR		NJ	NR
VALENCENE	UG/KG		NR	LN	R		NR		NR	NR		NR	NR.

		GS08SS		GS09SS		GS10SS		GS11SS		<b>GS12SS</b>		GS13SS		GS14SS
2 UNKNOWN COMPOUNDS	UG/KG	18	J		NR		ŇR	-	NR		NR		NR	NR
3 UNKNOWN COMPOUNDS	UG/KG	.,	NR		NR	29	7	35	J	30	7		NR	NR
ACETALDEHYDE	UG/KG		NR		NR		NR		NR		NR	16	NJ	12 NJ
BUTANAL	UG/KG	i i	NR		NR		NR		NR	5	Ŋ		NR	NR
HEXANAL	UG/KG	1	NR		NR	45	NJ	37	J	18	Ŋ		NR	NR
PENTANAL	UG/KG		NR		NR		NR		NR		NR		NR	NR
PROPANAL, 2-METHYL-	UG/KG		NR		NR		NR		NR		ŇR		NR	NR
UNKNOWN COMPOUND	UG/KG		NR		NR		NR		NR	-	NR	13	J	7 J
UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR		NA		NR		NR	NR

Table 10. Surface Soil Miscellaneous Analytical Results. Page 5 of 8

	Units	GS15SS	GS16SS	GS17SS	GS18SS	GS19SS	GS20SS	GS20SD	GS21SS
[3,4:9,10] DIBENZPYRENE	UG/KG	NR NR	NR	NR	NR	NR	NR		NR
1-DOCOSENE	UG/KG	NR	NR	NR	NR	NR	. NR	220 NJ	
1,2:3,4-DIBENZPYRENE	UG/KG	190 NJ	NA	290 NJ	NR	NR	NR	NR	NR
1,2:4,5-DIBENZPYRENE	UG/KG	! NR	NR	NR	NR	NR	NR	NR	NR
1,2:7,8-DIBENZPHENANTHRENE	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
1,2:7,8-DIBENZPHENANTHRRENE	UG/KG	NR NR	NR	NR	NR	NR	NR	. NR	NR
10 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR	NR	NR	NR	NR	NR	NR
10 UNKNOWN PAHS	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
11 UNKNOWN COMPOUNDS	. UG/KG	NR	NR	NR	NR	NR	NR.	NR	NR
11H-BENZO [A] FLUORENE	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
11H-BENZO [B] FLUORENE	UG/KG	350 NJ	NR	510 NJ	- NR	NR	NA	NR	NR
12 UNKNOWN COMPOUNDS	UG/KG	. NR	NR	NR	NR	NR	NR	NR	NR
14 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR	NR	NR	NR	NR NR	NR	3100 J
15 UNKNOWN COMPOUNDS	UG/KG	NR	NR	NA	NR.	NR	NR	NR	NR
16 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR	5900 J	NR.	NR	NR	NR	NR
17 UNKNOWN COMPOUNDS	UG/KG	NR NR	24000 J	NR	3100 J	NR NR	8800 J	9100 J	NR
19 UNKNOWN COMPOUNDS	UG/KG	12000 J	NB	NR	NR NR	NR	NR.	NR	NR
1-EICOSANOL	UG/KG	NR NR	NR	NR	NR.	NR.	190 NJ	NR	230 NJ
1-HEPTADECENE	UG/KG	NR.	NR	NR	NR.	NR	NR	NR NR	NR NR
1-OCTADECENE	UG/KG	NR	NR	NR	NR	NR	NR	NR NR	NR
1-TETRADECENE	UG/KG	NR.	NR	NR	NB	NB	NR NR	NR	NR NR
2 UNKNOWN KETONES	UG/KG	NR NR	NR.	NR	NR I	NR NR	NR.	NR	NR
2.2'-BINAPHTHALANE	UG/KG	NR	NR	NR	NR.	NR	NR	NR	NB
2.2'-BINAPHTHALENE	UG/KG	NR NR	NR	NR	NR NR	NR	NR.	NR NR	NR
21 UNKNOWN COMPOUNDS	UG/KG	NR NR	NB	NR	NA	NR	NR NR	NR NR	NR
25 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR	NR	NR	12000 J	NR	NR NR	NR
3 UNKNOWN COMPOUNDS	UG/KG	NR NR	NB	NR	NR NR	NR NR	NR.	NR NR	NR
3 UNKNOWN PAHS	UG/KG	NR	NR	970 J	NR NR	NR NR	NR	NR	NR NR
3.4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG	NR NR	NR	NR	NR NR	NR NR	NR	NR	NR.
4 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR	NR	NR NR	NR	NR	NR	NR
4 UNKNOWN PAHS	UG/KG	2000 J	NR NR	NR	NR NR	NR	NR NR	NR	NR
5 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR	NR	NR NR	NB NB	NR	NR	NR
5, 12-NAPHTHACENEDIONE	UG/KG	NR	NR NR	NR	NR NR	NR	NB	NR	NR
5.12-NAPHTAHCENEDIONE	UG/KG	NR NR	NR NR	NR	NR	NR.	NR.	NR	NR
5.12-NAPHTHACENEDIONE	UG/KG	NB	2600 J	800 NJ	NR NR	NR NR	NR	NR	NR NR
6 UNKNOWN COMPOUNDS	UG/KG	NR	NR	NR	NR	NB	NR	NR NR	NR NR
6 UNKNOWN PAHS	UG/KG	NR NR	NR NR	NR	NR NR	NR.	NR NR	NR NR	NR NR
7 UNKNOWN COMPOUNDS	UG/KG	NR	NR NR	NB	NR	NB	NR NR	NB	NR NR
7 UNKNOWN PAHS	UG/KG	NR	NA NA	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR
7H-BENZ [DE] ANTHRACEN-7-ONE	UG/KG	440 NJ	1600 J	390 NJ	NR NR	NR NR	NR NR	NR	NR NR
8 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR NR	NR	NR NR	NR NR	NR NR	NR NR	NR NR
8 UNKNOWN PAHS	UG/KG	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	- NR
9, 10-ANTHRACENEDIONE	UG/KG	NR	NR NR	NR	NR NR	NR NR	NR NR	NR NR	NR NR
9.10-ANTHRACENEDIONE	UG/KG	700 NJ	160 J	NR	NR NR	NR NR	NR NR	NR NR	NR NR
9H-FLUOREN-9-ONE	UG/KG	NR NR	NR NR	NR	NR NR	NR NR	NR NR	NR NR	NR
9-OCTADECENE, (E) -	UG/KG	NR NR	NR NR	NR	NR NR	NR NR	NR NR	NR NR	NR NR
ANTHRACENE, 2-METHYL-	UG/KG	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR
BENZ (A) ANTHRACENE-7, 12-DIONE	UG/KG	260 NJ	NR NR	NR NR	NR NR	NR I	NR NR	INR.	NR NR
BENZ (A) ANTHRACENE-7, 12-DIONE	UG/KG	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR
BENZ (A) ANTHRACENE, 7-METHYL-	UG/KG	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR
BENZ [E] ACEPHENANTHRYLENE	UG/KG	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR	NR NR
BENZO (B) NAPHTHO (2,1-D) THIOPH	UG/KG	NR NR	1400 J	NA NA	NA NA	NR NR	NR NR	NR NR	- NR
BENZO [B] NAPHTHO [2,1-D] THIOPH BENZO [B] NAPHTHO [2,1-D] THIOPHENE	UG/KG	NR NR	1400 J	310 NJ		NR NR			
BENZO [B] NAPHTHO [2,1-D] THIPHENE	UG/KG	NR NR	NA NA	NR NR	NR		NR NR	NR NR	NR
		NR NR			NR NR	NR	NR	NR	NR
BENZO [B] NAPHTHO [2,3-D] FURAN	UG/KG		NR NR	NR NR	NR NR	NR	NR NR	NR NR	NR NR
BENZO [8] TRIPHENYLENE	UG/KG	NR NR	NR NR	NR NR	NR	NR	NR NR	NR NR	NR NR
BENZO [E] PYRENE	UG/KG	NR	NR	880 NJ	NR NR	NR	NR	NR	NR

	Units	GS15SS	GS16SS	GS17SS	GS18SS	GS19SS	GS20SS	GS20SD	GS21SS
BENZO [J] FLUORANTHENE	UG/KG	i NR	NR	NR	NR	NR	NR	NR	NR
BENZOFLUORENE	UG/KG	NR NR	NR	NR	NR	NR	NR	NR	NR
CARYOPHYLLENE	UG/KG	. NR	NR	NR	NR	NR	NR	NR	NR
CHOLESTEROL	UG/KG	NR	NA	650 NJ	NR	. NR	NR	NR	NR
CHRYSENE, 1-METHYL-	UG/KG	300 NJ	NR	NR	NR	NR	NR	NR	NR
CHRYSENE, 5-MEHTHYL-	UG/KG	NR NR	NR	, NR	NR	NR	NR	NR	NR
CHRYSENE, 5-METHYL-	UG/KG	! NR	2300 J	310 NJ	NR	NR	NR	NR	NR
CHRYSENE, 6-METHYL-	UG/KG	NR	· NR	NR	NR	NR	NR	NR	NR
CYCLOPENTA (DEF) PHENANTHRENONE	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
D:C-FRIEDOOLEANAN-3-ONE	UG/KG	NR NR	NR	NR.	NR NR	NR	NR	NR	NR
HEXADECANOIC ACID	UG/KG	NR NR	NR	NR	110 NJ	160 J	130 NJ	150 NJ	NR
METHYLANTHRACENE	UG/KG	NR NR	NA	NR	NR	NR	NR	NR	NR
NAPHTHALENE, 2-PHENYL-	UG/KG	NR	NR	210 NJ	NR	NR	NR NR	NR NR	NR
PENTADECANOIC ACID	UG/KG	NR	NR	NR	NR	. NR	NR	NR	NR
PERYLENE	UG/KG	NR	200 J	NR	NR	NR	NR	NR NR	NR
PHENANTHRENE, 1-METHYL-	UG/KG	NR	NR	NR	NR	NR	NR	· NR	NR
PHENANTHRENE, 2,3-DIMETHYL-	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
PHENANTHRENE, 2-METHYL-	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
PHENANTHRENE, 4,5-DIMETHYL-	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
PYRENE 2-METHYL-	UG/KG	NR	NA	NR	NR	NR	NR	NR NR	NR
PYRENE, 1,3-DIMETHYL-	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
PYRENE, 1-METHYL-	UG/KG	NR	2000 J	NR	NR	NR	NR	NR	NR
PYRENE, 2-METHYL-	UG/KG	NR NR	1400 J	NR	NR	NR	NR	NR	NR
PYRENE, 4-METHYL-	UG/KG	NR NR	NR	NR	NR	NR	NR	NR	NR
STIGMAST-4-EN-3-ONE	UG/KG	NR	NR	NR	NR	NR	NR NR	320 NJ	NR
TESTOSTERONE	UG/KG	NR	NR	NR	NR	390 NJ	NR ·	NR	NR
TRIPHENYLENE, 2-METHYL-	UG/KG	NR	NR.	· NR	NR	NR	NR	NR	NR
UNKNOWN ALCOHOL	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
UNKNOWN ALKENE	UG/KG	NR	NR	NR	NR	NR	NR	NR NR	. NR
UNKNOWN CARBOXYLIC ACID	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
UNKNOWN COMPOUND	UG/KG	NR	NR	· NR	NR	NR	NR	NR	NR_
UNKNOWN PAH	UG/KG	NR	580 J	· NR	NR	NR	NR	NR	NR
VALENCENE	UG/KG	NR	1200 NJ	NR	NR	NR	NR	NR	· NR

	ſ	GS15SS	GS16SS	GS17SS	GS18SS	GS19SS	GS20SS	GS20SD	GS21SS
2 UNKNOWN COMPOUNDS	UG/KG	NR	16 J	NR	NR	NR	NR	. NR	NR
3 UNKNOWN COMPOUNDS	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
ACETALDEHYDE	UG/KG	24 NJ	NR	38 NJ	NR	18 NJ	NR	42 NJ	47 NJ
BUTANAL	UG/KG	NR	· NR	NR	NR	NR	NR	11 NJ	NR
HEXANAL	UG/KG	NR	31. NJ	NR	NR	NR	NR	NR	NR.
PENTANAL	UG/KG	NR NR	NR	NR	NR	NR	NR	NR	7 NJ
PROPANAL, 2-METHYL-	UG/KG	NR	NR	7 NJ	NR	NR	NR	7 NJ	NR
UNKNOWN COMPOUND	UG/KG	10 J	NR	12 J	NR	NR	NR	9 J	12 J
UNKNOWN COMPOUNDS	UG/KG	NR	NR	NR NR	NR	NR	NR	NR	NR

Table 10. Surface Soil Miscellaneous Analytical Results. Page 7 of 8

Miscellaneous Extractables						.,	
	Units	GS22SS		GS23SS		GS24SS	
[3,4:9,10] DIBENZPYRENE	UG/KG	:	NR		NR		NR.
1-DOCOSENE	UG/KG		NR		NR		NR
1,2:3,4-DIBENZPYRENE	UG/KG		NR		NR		NR
1,2:4,5-DIBENZPYRENE	UG/KG	1	NR		NR		R
1,2:7,8-DIBENZPHENANTHRENE	UG/KG		NR		NR		NR
1,2:7,8-DIBENZPHENANTHRRENE	UG/KG	3	NR		NR		NR
10 UNKNOWN COMPOUNDS	UG/KG	:	NR		NR		NR
10 UNKNOWN PAHS	UG/KG	1 .	NR		NR		NR
11 UNKNOWN COMPOUNDS	UG/KG	:	NR		NR		NR
11H-BENZO [A] FLUORENE	UG/KG	:	ÑЯ		NR		NR
11H-BENZO (B) FLUORENE	UG/KG	:	NR		NR		NR
12 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR
14 UNKNOWN COMPOUNDS	UG/KG	1	NR		NR		NR
15 UNKNOWN COMPOUNDS	UG/KG	7	NR	ĺ	NR	5900	J
16 UNKNOWN COMPOUNDS	UG/KG	•	NR		NR		NR
17 UNKNOWN COMPOUNDS	UG/KG	-	NR		NR		NR
19 UNKNOWN COMPOUNDS	UG/KG		NR	-	NR	·	NR
1-EICOSANOL	UG/KG		NR		NR		NR
1-HEPTADECENE	UG/KG	1	NR		NR	1 - 1	NR
1-OCTADECENE	UG/KG		NR		NR	<b>†</b>	NR
1-TETRADECENE	UG/KG		NR		NR		NR
2 UNKNOWN KETONES	UG/KG	:	NR	· · · · · · · · · · · · · · · · · · ·	NR	1	NR
2.2'-BINAPHTHALANE	UG/KG		NR		NR		NR
2,2'-BINAPHTHALENE	UG/KG		NR		NR	1	NR
21 UNKNOWN COMPOUNDS	UG/KG		NR	3600	J	1	NR
25 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR
3 UNKNOWN COMPOUNDS	UG/KG		NR		NR	1	NR
3 UNKNOWN PAHS	UG/KG		NA		NR	t	NR
3.4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG		NR		NR	<del>   </del>	NR
4 UNKNOWN COMPOUNDS	UG/KG		J		NR	<del>                                     </del>	NR
4 UNKNOWN PAHS	UG/KG		NR		NR		NR
5 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR
5. 12-NAPHTHACENEDIONE	UG/KG		NR	210	NJ	<del> </del>	NR
5,12-NAPHTAHCENEDIONE	UG/KG		NR	3.0	NR		NR
5.12-NAPHTHACENEDIONE	UG/KG		NR		NR	1	NR
6 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR
6 UNKNOWN PAHS	UG/KG		NR		NR	1700	
7 UNKNOWN COMPOUNDS	UG/KG		NR		NR	·	ŇR
7 UNKNOWN PAHS	UG/KG		NR		NR		NR
7H-BENZ [DE] ANTHRACEN-7-ONE	UG/KG		NR	210		140	
B UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR
8 UNKNOWN PAHS	UG/KG	·	NR		NR		NR
9, 10-ANTHRACENEDIONE	UG/KG		NR		NR		NR
9,10-ANTHRACENEDIONE	UG/KG		NR		NR -		NR
9H-FLUOREN-9-ONE	UG/KG		NR NR		NR		NR
9-OCTADECENE, (E) -	UG/KG		NR.		NR.		NR.
ANTHRACENE, 2-METHYL-	UG/KG		NR		NR		NR
BENZ (A) ANTHRACENE-7, 12-DIONE	UG/KG		NR		NR -		NR
BENZ (A) ANTHRACENE-7, 12-DIONE	UG/KG		NR NR		NR		NR
BENZ (A) ANTHRACENE, 7-METHYL-	UG/KG		NR	240			NR
BENZ [E] ACEPHENANTHRYLENE	UG/KG		NR	110			NR NR
	UG/KG		NR NR				
BENZO [B] NAPHTHO [2,1-D] THIOPH					NR		NR
BENZO [B] NAPHTHO [2,1-D] THIOPHENE	UG/KG		NR		NR		NR
BENZO [B] NAPHTHO [2,1-D] THIPHENE	UG/KG		NR I		NR		NR
BENZO [B] NAPHTHO [2,3-D] FURAN	UG/KG		NR		NR		NR
BENZO [B] TRIPHENYLENE	UG/KG		NR		NR		NR
BENZO [E] PYRENE	UG/KG		NR	240	NJ	840	NJ

	Units	GS22SS		GS23SS		GS24SS	1
BENZO [J] FLUORANTHENE	UG/KG	: "	NR	440	NJ		NR
BENZOFLUORENE	UG/KG		NR		NR		NR
CARYOPHYLLENE	UG/KG	:	NR		NR	1600	NJ
CHOLESTEROL	UG/KG		NR		NR		NR
CHRYSENE, 1-METHYL-	UG/KG	:	NR		NR		NR
CHRYSENE, 5-MEHTHYL-	UG/KG	7	NR		NR		NR
CHRYSENE, 5-METHYL-	UG/KG		NR		NR		NR
CHRYSENE, 6-METHYL-	UG/KG		NR		NR		NR
CYCLOPENTA (DEF) PHENANTHRENONE	UG/KG		NR	140	NJ		NR
D:C-FRIEDOOLEANAN-3-ONE	UG/KG	1	NR	•	NR		NR
HEXADECANOIC ACID	UG/KG	. 110			NR		NR
METHYLANTHRACENE	UG/KG		NR		NR		NR
NAPHTHALENE, 2-PHENYL-	UG/KG		NR		NR		NR
PENTADECANOIC ACID	UG/KG		NR		NR		NR
PERYLENE	UG/KG		NR		NE	260	
PHENANTHRENE, 1-METHYL-	UG/KG		NR		NR		NR
PHENANTHRENE, 2,3-DIMETHYL-	UG/KG		NR		NR		NR
PHENANTHRENE, 2-METHYL-	UG/KG		NR		NR		NR
PHENANTHRENE, 4,5-DIMETHYL-	UG/KG		NR		NR		NR
PYRENE 2-METHYL-	UG/KG		NR		NR		NR
PYRENE, 1,3-DIMETHYL-	UG/KG		NR		NR		NR
PYRENE, 1-METHYL-	UG/KG	!	NR		NR	. 160	NJ
PYRENE, 2-METHYL-	UG/KG		NR	130	2		NR
PYRENE, 4-METHYL-	UG/KG		NR		NR		NR
STIGMAST-4-EN-3-ONE	UG/KG		NR	170		2600	
TESTOSTERONE	UG/KG		NR		NR		NR
TRIPHENYLENE, 2-METHYL-	UG/KG		NR		NR		NR
UNKNOWN ALCOHOL	UG/KG		NR		NR		NR
UNKNOWN ALKENE			NR		NR	200	7
UNKNOWN CARBOXYLIC ACID	UG/KG		NR		NR		NR
UNKNOWN COMPOUND			NR		NR		ZR
UNKNOWN PAH	UG/KG		NR		NR		NR
VALENCENE	UG/KG		NR		NR		NR

#### Miscellaneous Volatiles

	Γ	GS22SS	• ]	<b>GS23SS</b>		<b>GS24SS</b>	
2 UNKNOWN COMPOUNDS	UG/KG	1	NR		NR	150	J
3 UNKNOWN COMPOUNDS	UG/KG		NR	30	J		NR
ACETALDEHYDE	UG/KG	1	NR.	14	NJ		NR
BUTANAL	UG/KG		NR		NR		NR
HEXANAL	UG/KG		NR		NR		NR
PENTANAL	UG/KG	!	NR		NR		NR
PROPANAL, 2-METHYL-	. UG/KG		NR		NR		NR
UNKNOWN COMPOUND	UG/KG	6	J		NR		NR
UNKNOWN COMPOUNDS	UG/KG	1	NR	,	NR	-	NR

#### **Data Qualifiers**

U-Analyte not detected at or above reporting limit. The number is the minimum quantitation limit.

J-identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-identification of analyte is acceptable; reported value may be blased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be blased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is 'average' of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | //-when no value is reported, see chiordane constituents | //-constituents or metabolites of technical chlordane

Table 11. Subsurface Soil Miscellaneous Analytical Results. Page 1 of 3

•	Units	GS01SB	GS02SB	GS03SB	GS06SB	GS07SB	GS08SB	GS09SB	GS10SB
1,2:3,4-DIBENZPYRENE	UG/KG	. NR	NR	NR	NR	NR	190 NJ	NR	NR
1,2:7,8-DIBENZPHENANTHRENE	UG/KG	NR NR	NR NR	NR	NR	NR	NR	NR	7400 NJ
1.2-BENZENEDICARBOXYLIC ACID	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
10 UNKNOWN COMPOUNDS	UG/KG	NA NA	NR	NR	NR	NR	. NR	NR	NR
11H-BENZO (B) FLUORENE	UG/KG	NR	NR	NR	NR	NR	NR	NR	9900 NJ
13 UNKNOWN PAHS	UG/KG	NR	. NR	NR	NR	NR	NR	NR	65000 J
1-PHENANTHRENECARBOXYLIC ACID	UG/KG	92 NJ	NA	NR	NR	NR NR	NA	NR	NR
2 UNKNOWN COMPOUNDS	UG/KG	NR	NR.	250 J	NR	NR	NR	NA	NR
2 UNKNOWN OCMPOUNDS	UG/KG	NR	220 J	NR	NR	NR	NR	NR	NR
2 UNKNOWN PHTHALATES	UG/KG	NR	NR	NR	NR	NR NR	NR	NR	NR
2,2'-BINAPHTHALENE	UG/KG	NR	NA	NR	NR	NR NR	NA	INR	3100 NJ
3 UNKNOWN COMPOUNDS	UG/KG	! NR	NR	NR	NR	NR	NR	370 J	RN
3 UNKNOWN PAHS	UG/KG	NR	NR_	NR	NR _	NR .	560 J	NR	NR
3,4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG	NR	NR _	NR	NR	NR NR	NR	NR	2800 NJ
4 UNKNOWN COMPOUNDS	UG/KG	NR.	NR_	NR	NR	NR	NR	NR	17000 J
5 UNKNOWN COMPOUNDS	UG/KG	840 J	NR	NR	NR I	NR.	NR	NR	NR
5,12-NAPHTACENEDIONE -	. UG/KG	NR.	NR _	NR	NR	NR	NR	NR NR	4600 NJ
5,12-NAPHTHACENEDIONE	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
7 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR	. NR	NR	NR.	1500 J	NR	NR
7 UNKNOWN PAHS	UG/KG	NR I	NR	NR	NR	_ NA	NR	NR	NR
7H-BENZ [DE] ANTHRACEN-7-ONE	UG/KG	NR.	NR_	NR	NR	NR	NR	NR	NR
7H-BENZ [DE] ANTHRACENEDIONE	UG/KG	NR.	NR	NR NR	NR	NR NR	NR	NR	4600 NJ
9,10-ANTHRACENEDIONE	UG/KG	; NR	NR _	NR NR	NR	NR	NR	NR	NR
BENZ [A] ANTHRACENE, 8-METHYL-	UG/KG	NR NR	NR	NR NR	NR.	NR.	NR	NR	2800 NJ
BENZO [B] NAPHTHO [1,2-D] THIOPHENE	UQ/KG	NR NR	NR	NR.	NR	NR	NR	NR	NR
BENZO [B] NAPHTHO [2,3-D] THIOPHENE	UG/KG	NR	NR	NR	NR	NR NR	NR	NR	4600 NJ
BENZO [B] TRIPHENYLENE	UG/KG	NR NR	NR	NR	NR	NR	NR	NR	NR
BENZO [E] PYRENE	UG/KG	i NR	NR _	NR NR	NR	NR	NR	NR	6200 NJ
CHRYSENE, 6-METHYL.	UG/KG	NR	NR	NR	NR	NR	NR	NA	6000 NJ
CYCLOPENTA (dEF) PHENANTHRENONE	UG/KG	NR	NR	NR	NR	. NR	NR	NR	NR
CYCLOPENTA (DEF) PHENANTHRENONE	UG/KG	NR	NR	NR	NR	NA	NR	NR	4700 NJ
ELLIPTICINE	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
NAPHTHO [1,2,3,4-DEF] CHRYSENE	UG/KG	NR	NR	NR	NR	NR	120 NJ	NR	NR
PERYLENE	UG/KG	NR NR	NR	NR	NR	NR	490 NJ	NR	16000 NJ
PYRENE, 1-METHY-	UG/KG	NR.	NR	NR	NR	NR NR	NR	NR	NR
PYRENE, 1-METHYL-	UG/KG	NR	NR NR	NR	NR NR	NR	NA	NR	9200 NJ
UNKNOWN AMIDE	UG/KG	NR NR	NR	NR	NR	NR	110 J	NR	NR
UNKNOWN CARBOXYLIC ACID	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR
UNKNOWN COMPOUND	UG/KG	NR NR	NR	NR	NR	190 J	NR	NR	NR
VALENCENE	UG/KG	NR	NR	NR	NR	NR	NR	NR	NR

		GS01SE		GS02SB	GS03SB		GS06SB	GS07SB		GS08SB	GS09SB	GS10SB	$\Box$
2 UNKNOWN COMPOUNDS	UG/KG		NA	NR	I N	Ä	NR		NR	NR	RN	25 J	$\Box$
UNKNOWN COMPOUND	UG/KG		NR	NR	N	R	8 J		NA	8 J	NA	NF.	a ]

Table 11. Subsurface Soil Miscellaneous Analytical Results. Page 2 of 3

	Units	GS11SB		GS13SB		GS14SB		GS15SB		GS16SB	GS18SB	1	GS19SB
1.2:3,4-DIBENZPYRENE	UG/KG	,	NR		NR		NR	1	NR	NR		NR	NR NR
1.2:7.8-DIBENZPHENANTHRENE	UG/KG		NR		NR		NR	1	NR	NR.		NA	NR
1.2-BENZENEDICARBOXYLIC ACID	UG/KG		NR		NR		NR	1	NR	NR.	120	NJ	NR
10 UNKNOWN COMPOUNDS	UG/KG		NR		NR	3900	J	1	NR	N8		NR	NR NR
11H-BENZO [B] FLUORENE	UQ/KG		NR		NR		NR	T	NR	NR		NR	NR NR
13 UNKNOWN PAHS	UG/KG		NR		NA		NR		NR	NR		NR	NR
1-PHENANTHRENECARBOXYLIC ACID	UG/KG	-	NR		NR		NR	1	NR	NR.		NR	NR.
2 UNKNOWN COMPOUNDS	UG/KG		NR	1000	J		NA		NR .	220 J	450	7	NA NA
2 UNKNOWN OCMPOUNDS	UG/KG		NR		NR		NR		NR	NR		NR	NR NR
2 UNKNOWN PHTHALATES	UG/KG		NR		NR		NR		NR	NA	230	5	· NR
2,2'-BINAPHTHALENE	UG/KG	i	NR		NR		NR	1	NR	NR		NR	NR
3 UNKNOWN COMPOUNDS	UG/KG	530	J		NR -		NR		NA	NR		NR	NR
3 UNKNOWN PAHS	UG/KG		NR		NR		NR		NR	NR		NA	NR
3,4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG		NR		NR	290	NJ		NR	NR		NR	NR NR
4 UNKNOWN COMPOUNDS	UG/KG	:	NR		NR		NR	1	NR	NR		NR I	NR
5 UNKNOWN COMPOUNDS	UG/KG	1	NR		NR		NR	<del></del>	NR	NR	1	NR	NR
5,12-NAPHTACENEDIONE	UG/KG	1	NR		NR		NR	1	NR	NR		NR	NR NR
5,12-NAPHTHACENEDIONE	UG/KG		NR		NR	290	NJ	1	NR	NR		NR	NR
7 UNKNOWN COMPOUNDS	UG/KG		NR		NR		NR		NR	NR.		NR	NR NR
7 UNKNOWN PAHS	UG/KG	1	NR	1	NR	3200	5	1	NR	NR		NR	NR
7H-BENZ (DE) ANTHRACEN-7-ONE	UG/KG		NR		NR	500	NJ	1	NR	NR	1	NR	NR
7H-BENZ [DE] ANTHRACENEDIONE	UG/KG	1	NR		NR		NR	1	NR	NA		NR	NR
9,10-ANTHRACENEDIONE	UG/KG		NA		NR	320	Z		NA	NR		NR	. NR
BENZ (A) ANTHRACENE, 8-METHYL-	UG/KG		NR		NR		NR		NR	NR		NR	NR
BENZO [B] NAPHTHO [1,2-D] THIOPHENE	UQ/KG		NR		NR	270	NJ		NR	NR		NR	NR
BENZO (B) NAPHTHO [2,3-D] THIOPHENE	UG/KG		NR		NR		NR		NR	NR		NR	NR
BENZO (B) TRIPHENYLENE	UG/KG		NR		NR	340		I	NR	NR		NR	NR
BENZO [E] PYRENE	UG/KG		NR		NR	1900	NJ		NA	NR		NR	NR
CHRYSENE, 5-METHYL-	UG/KG	1	NR		NR		NR		NR	NR		NR	NR
CYCLOPENTA (dEF) PHENANTHRENONE	UG/KG		NR		NR	320			NR	NA		NR	NA NA
CYCLOPENTA (DEF) PHENANTHRENONE	UG/KG		NR		NR		NR		NR	NR		NR	NR
ELLIPTICINE	UG/KG		NR		NR	270			NR	NR		NR	. NR
NAPHTHO [1,2,3,4-DEF] CHRYSENE	UG/KG		NR		35		NR		NR	NR		NR	NR
PERYLENE	UG/KG		NR		5	420			NR	NR		NR	NR
PYRENE, 1-METHY-	navka		NR		35	340			NR	NR		NR	NR
PYRENE, 1-METHYL-	UG/KG		NR		NA		NŘ		NA	NR		NR	NR NR
UNKNOWN AMIDE	UQ/KG		NR		VR.		NR		NR	NR		NR	NR
UNKNOWN CARBOXYLIC ACID	UG/KG	86			VR.		NR		NR_	NR		NR	NR
UNKNOWN COMPOUND	UG/KG		NR		NR		NR	310		NA		NR	95 J
VALENCENE	UG/KG		NR		VR.	1	NR		NR	NR		NR	NR

	GS11SB	GS13SB	GS14SB	GS15SB	GS16SB	GS18SB	GS19SB
2 UNKNOWN COMPOUNDS UG/K	g NR	NR	NR NR	20 J	NR	NR	NR
UNKNOWN COMPOUND UG/K	G NR	NR.	8 J	NR	NR	NR	NR

Table 11. Subsurface Soil Miscellaneous Analytical Results. Page 3 of 3

•	Unite	GS20SB	l I	GS21SB	
1,2:3,4-DIBENZPYRENE	UG/KG		NR		NR
1,2:7,8-DIBENZPHENANTHRENE	UG/KG		NR		NR
1,2-BENZENEDICARBOXYLIC ACID	UG/KG		NR		NR
10 UNKNOWN COMPOUNDS	UG/KG		NR		NR
11H-BENZO (B) FLUORENE	UG/KG		NR		NR
13 UNKNOWN PAHS	UG/KG		NR		NR ·
1-PHENANTHRENECARBOXYLIC ACID	UG/KG		NR		NR
2 UNKNOWN COMPOUNDS	UG/KG		NR		NR
2 UNKNOWN OCMPOUNDS	UG/KG		NR		NR
2 UNKNOWN PHTHALATES	UG/KG		NR		NR
2,2'-BINAPHTHALENE	UG/KG		NR		NR
3 UNKNOWN COMPOUNDS	UG/KG		NR		NR
3 UNKNOWN PAHS	UG/KG		NR		NR
3,4-DIHYDROCYCLOPENTA (CD) PYRENE	UG/KG		NR		NR
4 UNKNOWN COMPOUNDS	UG/KG		NR		NR
5 UNKNOWN COMPOUNDS	UG/KG :		NF		NR
5,12-NAPHTACENEDIONE	- UG/KG		NA		NR
5,12-NAPHTHACENEDIONE	UG/KG		NR		NR
7 UNKNOWN COMPOUNDS	UG/KG		NR		NR
7 UNKNOWN PAHS	UG/KG		NA		NR
7H-BENZ [DE] ANTHRACEN-7-ONE	UG/KG		NR		NR
7H-BENZ [DE] ANTHRACENEDIONE	UG/KG		NR		NR
9,10-ANTHRACENEDIONE	UG/KG		NR.		NŘ
BENZ [A] ANTHRACENE, 8-METHYL-	UG/KG		NR		NR
BENZO [B] NAPHTHO [1,2-D] THIOPHENE	UG/KG		NR		NR
BENZO [B] NAPHTHO [2,3-D] THIOPHENE	UG/KG		NR		NR
BENZO (B) TRIPHENYLENE	UG/KG		NR		NR
BENZO [E] PYRENE	UG/KG		NR		NR
CHRYSENE, 5-METHYL-	UG/KG		NR		NR
CYCLOPENTA (dEF) PHENANTHRENONE	UG/KG		NR		NR
CYCLOPENTA (DEF) PHENANTHRENONE	UG/KG .		NR		NR
ELLIPTICINE	UG/KG		NR ]		NR
NAPHTHO [1,2,3,4-DEF] CHRYSENE	UG/KG		NR		NR
PERYLENE	UG/KG		NR		NA
PYRENE, 1-METHY-	UG/KG		NR		NR
PYRENE, 1-METHYL-	UG/KG		NR		NR
UNKNOWN AMIDE	UG/KG		NR		NR
UNKNOWN CARBOXYLIC ACID	UG/KG		NR		NR
UNKNOWN COMPOUND	UG/KG	130		84	Ĵ
VALENCENE	UG/KG	120	NJ		NR

#### Miscellaneous Volatiles

	<u> </u>	GS20SB	GS21SB
2 UNKNOWN COMPOUNDS	UG/KG	NR NR	NR
UNKNOWN COMPOUND	UG/KG	: NR	NR

#### **Data Qualiflers**

U-Analyte not detected at or above reporting limit. The number is the minimum quantitation limit. U-identification of analyte is acceptable; reported value is an estimate.

UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate.

N-Presumptive evidence analyte is present; analyte reported as tentative identification.

NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

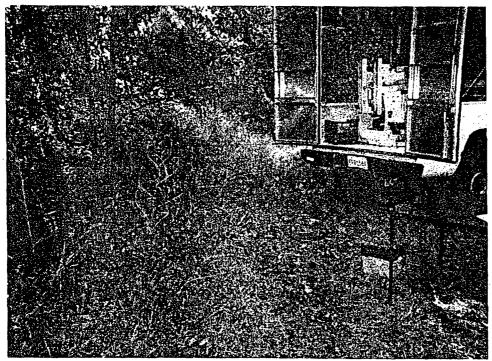
K-identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-identification of analyte is acceptable; reported value may be bigsed low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to interferences. | A-Analyte analyzed in replicate. Reported value is 'everage' of replicates.
R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

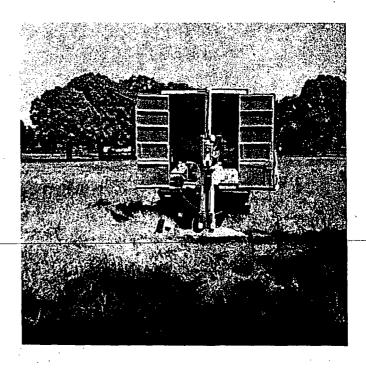
Appendix C Site Photographs

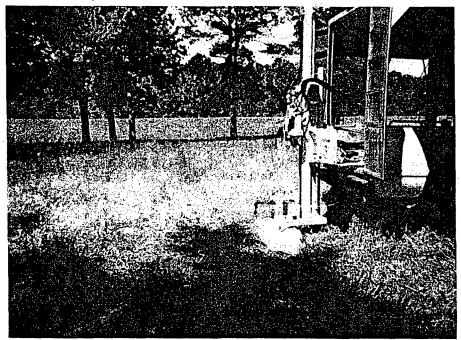
# Groundwater



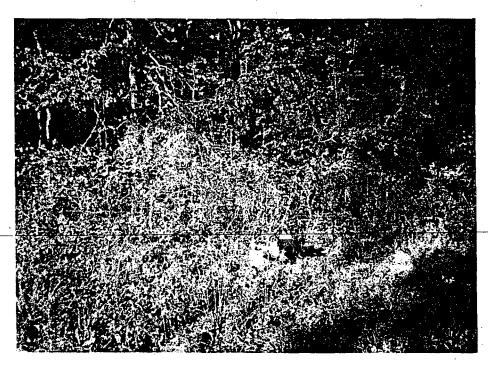
Site of GS-01-SS,SB,GW (Roll 3, frame 24)

Site of GS-02-GW,SS,SB, facing east (Roll 3, frame 21)

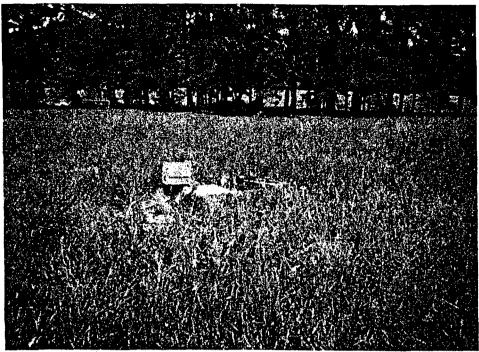




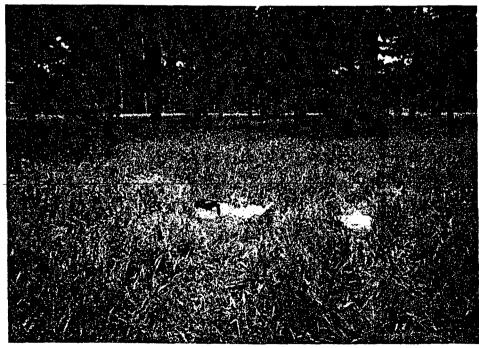
Site of GS-02-GW, looking north (Roll 3, frame 22)



Site of GS-03-GW, edge of woods surrounding Creosote Slough (Roll 3, frame 20)



Site of GS-04-SS,SB,GW, facing west (Roll 1, frame 19)

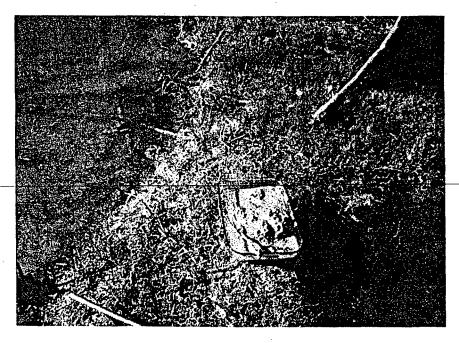


Site of GS-05-GW/SS/SB, facing west (Roll 3, frame 18)

# **Sediment**



Site of GS-01-SD (Roll 2, frame 23)



Sample GS-01-SD in pan (Roll 2, frame 22)



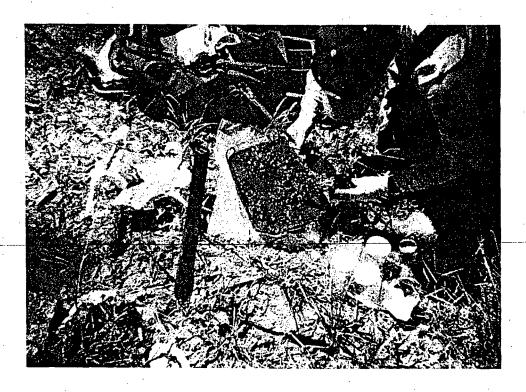
Site of GS-02-SD, facing east (Roll 2, frame 21)



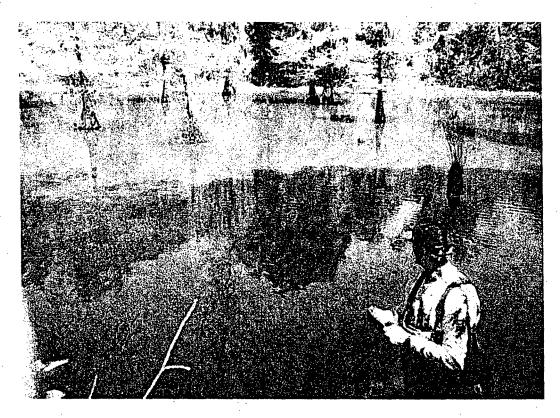
Sample GS-02-SD in pan (Roll 2, frame 20)



Site of GS-03-SD, facing east (Roll 2, frame 19)



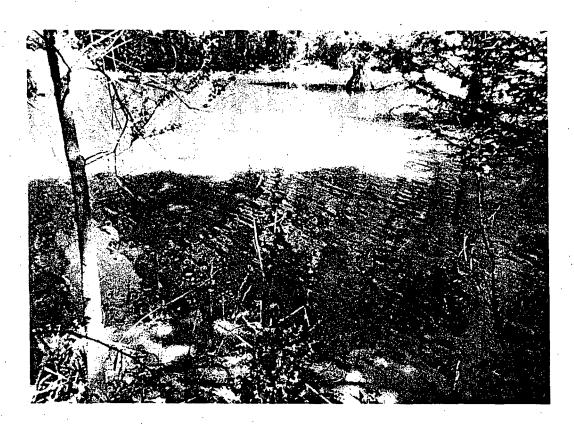
Sample GS-03-SD in pan (Roll 2, frame 18)



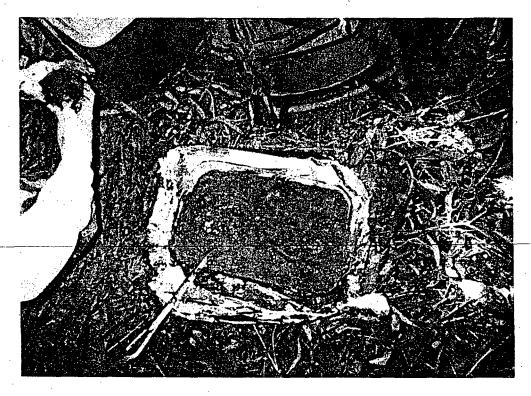
Site of GS-04-SD, facing north (Roll 2, frame 17)



Sample GS-04-SD in pan (Roll 2, frame 16)



Site of GS-05-SD (Roll 2, frame 15



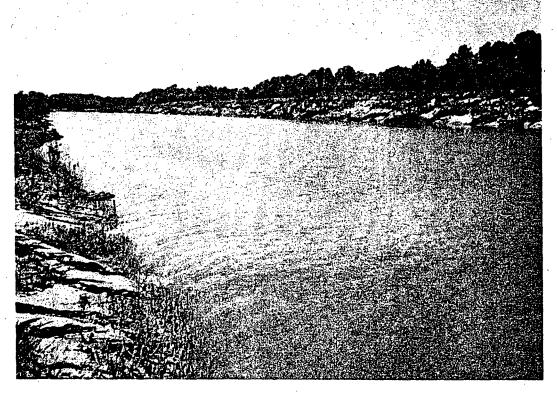
Sample GS-05SD in pan (Roll 2, frame 14)



Site of GS-06-SD (Roll 2, frame 11)



Sample GS-06-SD in pan (Roll 2, frame 10)



Site of GS-07-SD (Roll 2, frame 13)

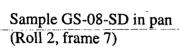


Sample GS-07-SD in pan (Roll 2, frame 12)



Site of GS-08-SD (Roll 2, frame 8)

Site of GS-08-SD (Roll 2, frame 8)

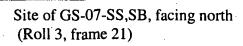




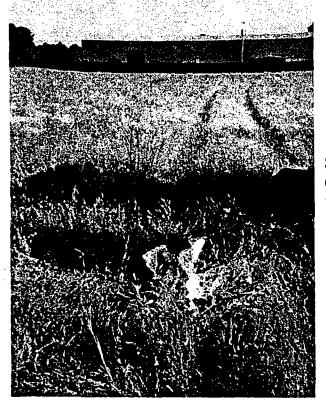
# Soil Sample Locations



Site of GS-06-GW,SS,SB, facing north (Roll 3, frame 24)

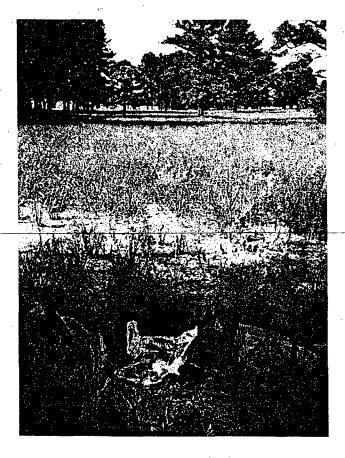






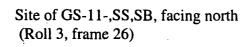
Site of GS-08-,SS,SB, facing north (Roll 3, frame 22)

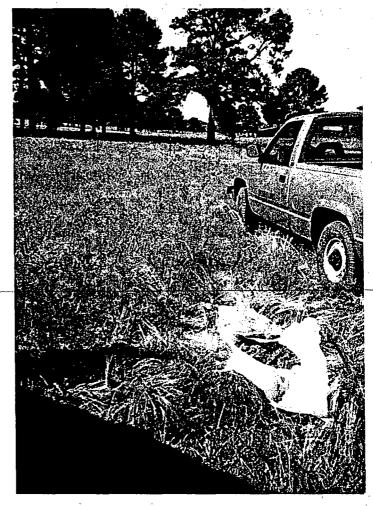
Site of GS-09-,SS,SB, facing north (Roll 3, frame 25)





Site of GS-10-,SS,SB, facing north (Roll 3, frame 27)

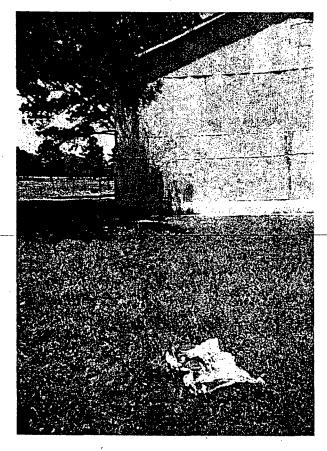






Site of GS-12-,SS,SB, facing north (Roll 3, frame 23)

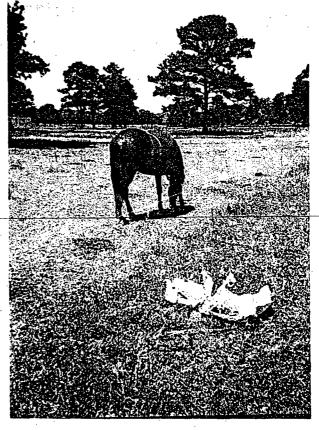
Site of GS-13-,SS,SB, facing north (Roll 3, frame 31)





Site of GS-14-,SS,SB, facing north (Roll 3, frame 28)

Site of GS-15-,SS,SB, facing north (Roll 3, frame 30)





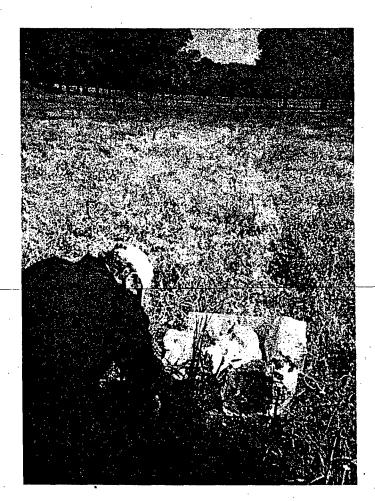
Site of GS-16-,SS,SB, facing north (Roll 3, frame 33)

Site of GS-17-,SS,SB, facing north (Roll 3, frame 35)





Site of GS-18-,SS,SB, facing north (Roll 3, frame 37)



Site of GS-19-,SS,SB, facing north (Roll 3, frame 36)



Site of GS-20-,SS,SB, facing north (Roll 3, frame 34)

Site of GS-21-,SS,SB, facing north (Roll 3, frame 32)





Site of GS-23-SS, facing north (Roll 3, frame 29)

### Appendix D Log Book Photocopies

7-22-23

FLD WOOD MS

CLEAS WASD

THOMPA

AND METALS

- WILL COLLECT 3 ENCORE

3- SON WHICH INCCUPED SPLITTER OWNER

DEPTH 0-6 -55 24-36 -5B

AT EACH LOCATION WOTH IS A MEEL UNIES OFFERWER NOTED.

USED DEET

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65-18-55 N 3611887, US	G5-19-55	N 36123	
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		GARY/BAANGE C	Ay	
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1355 TIONE 1315 0-6" FARRIAN COLCACITA THOM ON MISTO EXILD B. 501 BACKIN SOLL 3-33 13 SB TAME 1325 24-38" FAMILE COLLEGATE PARMAN DILLEN/nitto 3-6 NC ORANDE CLAY (SANDY) 3 dra PICTURE 7 LADICING NOATH

MYE 1340
FARRIER COLECTED
THOMAN MITER/FILLED 3-ENC
1-29
3-80

PARRIER COLLECTED

TOWN AND MIXED FILLED 3 ENC

1-23

BROWN CRAY

PICTURE & CORNE NORFOL

4-22/0 0-6 09-55 11 - 55DONE 1615 TIME FARRIER COLLECTE THOMAN CONFERDAJNIKED/814BS 3-ENC 1600 BROWN Snally Jox 11-58 24-36 08-58 MME1625 TIME FARMIEN COLLEGIED THOMON MINED /FILLD 3-ONC BROWN CLAY PICTURE 12 LOUGHT NORTH

0-6 FARRIES (DICEOSE THOMAN MIXED/BICKED 3-ENC 1-29 7-89. BROWN CLAY-WER

FARRIER COLLECTED THOMAN MIXED STIKED

3-ENC

BNOWN CLAR

PRTURE 13 LOOKING NORGET

	ILLO SAFERY METERNY DNSIFE OZIN
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DHUMAN MIXED FILLED 3-ENC	FARRIZA COXLECTIO
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	3-89
	BRUUN IPMPT 53/L
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6-SB 24.36" DME 650	12-58 24-36
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	GRPY CIAY
PICTUBE 14 LOWENS NOTTH	
	PICTURE 15 LOOKING NORTH .

\*

4-23 000 05-55 0-6' DYK 09/5 TRABIES CONSCIONED THOMAN MIXED/FILEDD 3-400 BRUNN SANDY SOLS 05-5B 2K-36" PINE 0925 FAMILY COLLEGED DAMAN MITTER FILLS GARYCLARY - WIET PICKYNE 18 LOOKING NORTH

1625 Flowood Drive Flowood, MS

USEPA 980 College Station Rd 1-11/445, GA 30605

Brian Striggen - project lander

Week of April 20, 2003.

Bubby Lewis

J woll lov i

Wook 5

Tim Simpson

Start Time 7:00

- SHELY muling held. Tim Simpson is SSO.

Simpson / Lewis ... Sortiment Sampling Thomas / Farrier -

Simples Allen - ground water

Strugon - Project Leader

Sile Reion done on 4-20-2003 by striggow, Lewis,
Simpson + Frence.

Departed Athens via GOU to sile on 4/21/03

- Svony clear , waren

$\overline{(2)}$		·	(j)
		· ! .	V/>>/
GSOISD Background Sample			1/22/03 Canadian Goese, crow
	L Cricker, Frency, Gerr	ph picorei	
Sampler BLOWIS Lordomer sent longer by Sampson	Sample chain of a		maintained  g 9 Stanles's Steel
Samples callected - TOC	Scoop, spoon	+ 9 Ta	or por VOCS
- VOLS (NOT mixed), % Moisture -SNOCS/POST/PAH -metals	Sompler		VIE AL LUCORE
sample - Fine clay -gray -coarse organic matter	CAMPLE LOENTIAN		
- pine st-aw + sticks Removed from		<u> </u>	
Sample		oAth!	
Photos 1+Z - Large Cottonmorth - B Lowis, 904	7/26		trac
Photo 3 - Background Location in Bayon  905, Lewis, Facing SW (2300)			
photo 4 - sample in pan			
Photo 4 - Sample in part 905, Levis			co Background well
gps data collected By B. Lewis			
Location - Jockson Prep Bayon Left of 'porth way"			
Lett of just hong	ے. ۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔۔	4001	
Sander reed in cooler			

GS02 SD sampler - Blowis Sample time - 950 Samples - TOC, metals, SVOC/POST/PAH, To moisture stunder steel sampling equipment + gloss pon used sample - gray I brown - siltly clay mix some organic material Boot mass Removed, Approx. 4 doep where sample collected slight sheen seen in sample pan - photo 5 - B lows 955 . Tring West, Nur 310" photo 6- B Lewis 955 Sample 02 BAYOU DZ location is SW of sile - Location about & A from Bent 195 coordinate collected by Bleuis Approx 20 ft From drainage NW Bayou Moad way (slag) 5W BAY OU

[6503 SD	4/22/03
1 sampler B LOWIS	
collected - TCC, SUDC, June stare, VOC., Pesil /PCB/PAH	
Stainless steel scoop, spoon + glass p. boundgray sonnie - heavy wlorg An material word mass	ic (corre
Note-shown seen. Creosole sine detected	
Photo 7 - Sample location 945 - Facing East 130°	
photo 8 - Sample 03; lenis, 945	8 
grs dula collected by B. Lewis	
Location From non vegetative 1 Emergent Vegetier neur bank	voc(

650750

Sampler - B. Lewis Surple 1245

Camples - TOE, UK, 90 mostro, metals

Location - Pent Pour downstreams

Sample was offer Sand

phits 13 Location - Lewis 1248

Stamless Steel scaplspoon used Mixed in glass pan.

glass 8-02 juss used for All Samples except To moisture (24/21 (2 02 glass) + Encoves (for UOLS)

GS 0830 Provice Branch

Samples - B. Lives samples TOC, To misting, VOCs, SUOC., met

photo 18 Location 19 Sample

Sediment was fine brown sand Prairie (reak ) 1st had sandx Shoots along ide length no organic much was excontered

VOCS : directly in to Encores (offerted at Stainless sheet scoop. mixed in glass pan

> Bardid water spate spring in Crick

4/27/2003 1122/03 680 GS24SS 16506 Sp Penal Buer location, upstream of Prairie Branch 0-6" surface soil samples
collected without sugar is steel)
mixed in gloss pan wis steel spoon Collieted Sangle 1515 Collected 3 Encore & 3 802 g/g55 E/M/Tax - photo 15 Cocation 1523 - photo 16 Sample 1529 Sampler sample time -16 35 collected. netals, SUOCs/pesT/FAH, To moisture collected 2 Extra 8 02 to glit Sedment use fine to medion The whole river was sandy

no organic much observed phado 22 Lucation by Linis 23 Sample by lines Soil was donk brown organic rich Photo 17, Lowis, @ 1530 -Luc mussels in shallows along River Mosquis fish Collectin 2 6" Augre side by side Ideal fied . Other large Fish seen hitting surface collected wisisted Augur
mixed in glass pan wisisted spoon VOCS - directly into Encore Low in 1819)

G503 4/23/12 Surface Soil 55. Collection Time - 740 By, Blows Samples collected - VOL (4 Enlares), To mudace MSINSD - SUCCEPPANFEST, met - Extra 8 02 tor split Brown I gray (W/some orange) clay. Some organic material (Black natural) present gos data by bonty Lewis Subs-- love 501 sampler - B. Lowis depin - 12-18", due la water we sampled at 12.15" Sample Fine - 755 collected - Vol, SVIX (PAH /POST Mutals, go moistire Extra 802 for split Brown clay location - near will 3 on edge of photo 24 by lows @ 812 Francy south Sample location 03

Surface Soil - 0-6"

Surplus Blows

Sumple time - 0830

collected - bot. To mosture, Mutals, SVOC PES PAH

sisc collected an extra 302 for spills

gray brown clay

Sumple time 840

Sampler - Simpson

sticky gray (brown clay, some crange

collected - SUOC/PAILIPEST, met, VOC, 70 meisture, cithra 802
photo 25, canky a 2 @ 885 by Lenis

Of location, facing enst.

GPS dada collected by Bobby lewis
surpled wisternless steel hand Ruger

03-0474

Book 4

Crew: Simmons/ Allen

## Photograph Log Comen #3

- 1. 65-01-55/58 locations
- 2. 65-01-55/58 locations w/ Graphotes in bedrach
- 3. 65-01-6W Well Pringing)
- 4. 65-02-55/53/GW lasking North
- 5. In lauking East
- 6.65-03-6W looking West-
- 7.65-04-6W looking West
- & 65-05-6W lasking Nest
- 9. Geogrape rad extraction @ Southard ofte
- 10. Geophobe not extraction & sontand site

C530	Arrived at Indeson Preparations School
	property for bordegiound sample G5-01
	Collected 01-55 @ 0'-6"
*	n Energy (ms/mso)
	252
	Brownish red sondy clay
0915	Collected GS-01-5B @ 24"-36"
	3 Encores 1 2 oz
	2 802
	Brownish red soiled clay, more most
	Location moved 12" due to obstruction.
	rosultar word 15 and 12 absurgilar.

2	
H20 Instrument calibration	65-01-GW well put in by Geoprobe
Nech 2100 Turbithination #9 5td values  low std = 4.83  4.78	Total well depoth 200 Screen @ 16-20
$\frac{1000 \text{ std}}{1000 \text{ std}} = 4.83$ 4.78 $\frac{1}{56.4}$	Death to NEO 5.6 (from grand level)
high std = 532 530	
(	1000 Bergin parge of well
Orion. 235A pl #6	
Self trest - OK	Time of Condus NTLL
7 st = 7.38 us	1503 54+ 186 251 -
4 std = 4.01	1010 5,51 19.3 214 344
$ s _{\infty} =  s _{\infty} =  s _{\infty}$	1016 5.58 19.2 214 42.8 1026 5.66 18.2 213 13.3
Drisi 140 Conductivity # 710151	1032 5.70 18,2 215 8,53
46.7 = 46.7	1539 - 6.24
445 = 427	1042 Began allecting metals, but turbulity
<u> </u>	went buck up to 7 12 Met Continues
	1051 - 187
	1062 184 219 300
	1057 6.69
	100 5,41 H.5 218 5.97
	1105 Collected water sample
· · · · · · · · · · · · · · · · · · ·	& IL ember (MS/MSD)
	$\leq 16$ poly in
	Sompled motals that pamp.
	- wither wait? Will brille.

65-02-55

1313 Collected sample from 0'-6".

3 Encores

1 202 3 802 (1 for property owner) Brown organic suit, flood plain

1334 Gilladed G5-02-53 From 24'-36"

3 En6125

3, 802 (1, for owner)

Ton clay soil

1350 Began well installation for GS - OZ - GW Total depth 20 0 Screen & 16'-20'

1408 Brogn pirale

 Time,	ph	02	Cond	NIL
 1410	5.35	19.1	750	
 1420	4.89	18.4	726	135
1430	4.91	18 4	.722	17.2
1440	454	18.5	726	5.97
 1450	495	18.4	724	3.70
		-1.		

1465 Simpled Well GS-02-CW Turbialty increased, purged note for metals sample

1642 - Collected 65-22-55 from 0'-6" Ton sonly clay

	i .				•	
4/23/03 G5-03-6W Jocation						
0735 Instrument contribution	0740		Goopraly	100 of C	35-03-0	GN
	   (					
12 21008 Turbidinatur # 9 St. velus	<u> </u>	total douth	, 20	Scrazn (	et 16-20	
) ow std = 485 478		Dooth to He		below of	send surfax	<u>C</u>
mid stel = 56.7 56 A	0530	Stirted per	je.	·		
high std = 532 530	, , , , , , , , , , , , , , , , , , ,	· · · · · · · · · · · · · · · · · · ·		· · · · · · · · · · · · · · · · · · ·		
	Tine	PH		(3nd 1/15		
Orion 230M all #6	2020	651	17.7	x377 40	5 —	instal
7 sid = 7.04	0507				<u>E3</u>	
4 std = 4.00	38/2	5 59	175	186	7)333	
s) spi = 102 6	0522	5.92	174	181	274	
10 Ad chade = 1013	0835	5.93	17.5	181	126	
Orion 110 Conductivity #710151	0542	5.93	17.4	120	7)	
46.7 = 47.0	_ 0%52				55.5	
445 = 431	V657		<u>!</u>		49.0	
	0907				31.9	
	0722		17.5	180	19:7	
1600 Pet collection check	9732				14.4	
Hach 2100P #9	0942		<u>.</u>	· · · · · · · · · · · · · · · · · · ·	12.2	
lay std = 4.79	0947		<u></u>	·	10.2	
md std = 56.3	0752		<u></u>		984	· '. ====================================
high 57d = 528	· 		· · · · · · · · · · · · · · · · · · ·	<b></b>		
	<u> </u>	Silvated	Sony 2	(5-03-6)	<b>V</b>	
Or,50 230 A & #6	· · · · · · · · · · · · · · · · · · ·					
4 = 44			: : i			
1 = 7.2			<del></del>			
	<del></del>			· · · · · · · · · · · · · · · · · · ·		
Orran Mi Conductività	÷		İ	-		

65-04-6W Ystal depth 20 Screen @ 16-20! G5-05-6W Total depth

σc[j

Bearn Puroje of GS-134-6W 104-5 Time 18.2 1545 intal 60 15.4 9.84 Bes 17.9 11:10 5.86 11.25 5.91 18.2 58 15.2 50) 14.6 11:40 5.99 18.5 61 12:30 191 5.68 10.7 12:45 6.00 60 19.0 11.7 5.77 19.0 11.0 13:00 11.0 199 13:15 5.81

Collected sample GS-04-GW

1320

Began arge & GS-05-GW off Time Cond initial 1105 696 11:12 5.39 17/6 595 376 1122 5.43 17:6 603 16.1 1132 5.53 176 605 8,60 1142 17/6 5.32 807 6.23 alected sande 1)45 05-GW

Appendix E Laboratory Data Sheets

# USEPA Contract Laboratory Program Organic Traffic Report & Chain of Custody Record

Case No:

DAS No:

31635

	<u> </u>						· · · · · · · · · · · · · · · · · · ·	
Region:	4		Date Shipped:	4/22/03	Chain of Custody Re	ecord	Sampler Signature:	
Project Code:	03-0474	]	Carrier Name:	FedEx	.]		Signature.	
Account Code:	50102D04ZZQB00		Airbiil:	836362682006	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:	İ	į	Shipped to:	Liberty Analytical	1			
Spill!ID:			omphou to.	Corporation				
Site Name/State:	Gulf States Creosoting/MS	1		501 Madison Avenue	2		·	
Project Leader:	Brian Striggow	l		Cary NC 27513	3			
Action:	Screening Site Investigation	1		(919) 379-4100				
Sampling Co:	USEPA SESD				4			

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION	SAMPLE COL DATE/TIM		INORGANIC SAMPLE No.		QC Type
D1XR4	Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-01-SD	S: 4/22/03 9:	00	MD1XR4		
D1XR5	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-18-SB	S: 4/22/03 8:	50	MD1XR5		
D1XR6	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-18-SS	S: 4/22/03 8:	40	MD1XR6	(	
D1XR7	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-19-SB	S: 4/22/03 9:	35	MD1XR7		
D1XR8	Surface Soil (0"-6")/	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-19-SS	S: 4/22/03 9:	30	MD1XR8		· · · ·
D1XR9	Dan Thoman Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-02-SD	S: 4/22/03 9:	50	MD1XR9		
D1XS0	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-17-SS	S: 4/22/03 10	05	MD1XS0		
D1XS1	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-17-SB	S: 4/22/03 10:	20 .	MD1XS1		
D1XS2	Surface Soil (0"-6")/	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-20-SS	S: 4/22/03 10	45	MD1XS2		
D1XS3	Dan Thoman Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-20-SB	S: 4/22/03 10	55	MD1XS3		

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	D1XS7, D1XT2	Rolf De O Many	
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?
MOICA = CLP %Moistu	re/Calcareous (soil), VOA-s = CLP TCL Volatiles (soil)		

TR Number: TR Number: 4-252433676-042203-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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# USEPA Contract Laboratory Program Organic Traffic Report & Chain of Custody Record

Case No:

DAS No:

31635

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Region: Project Code:	4	1	Date Shipped:	4/22/03	Chain of Custody	Chain of Custody Record		Sampler Signature:		
Account Code:	03-0474 50102D04ZZQB00	li i	rier Name:	FedEx	Relinguished By	(Date / Time)	Received By	(Date / Time)		
	90102D0422GB00	Alrbill:	oill;	836362682006	Reiniquished by	(Date / Inne)	Received by	(Date / Time)		
CERCLIS ID:		Shin	pped to:	Liberty Analytical	1		1			
Spill ID:		1	spou to.	Corporation						
Site Name/State:	Gulf States Creosoting/MS			501 Madison Avenue Cary NC 27513 (919) 379-4100	2					
Project Leader:	Brian Striggow	1			2			•		
Action:	Screening Site Investigation				3	· · · · · · · · · · · · · · · · · · ·				
Sampling Co:	USEPA SESD				4 .					

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		COLLECT E/TIME	INORGANIC SAMPLE No.	QC Type	
D1XS4	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-20-SD	S: 4/22/03	10:45	MD1XS4		
D1XS5	Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-03-SD	S: 4/22/03	10:30	MD1XS5		
D1XS6	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-16-SS	S: 4/22/03	11:30	MD1XS6		
D1XS7	Subsurface Soil (24"-36")/ Dan Thoman	IJĠ	MOICA (21), VOA-s (21)	(Ice Only) (10)	GS-16-SB	S: 4/22/03	11:40	MD1XS7	<del></del>	
D1XS8	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-21-SS	S: 4/22/03	12:05	MD1XS8	<b></b>	
D1XS9	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-21-SB	S: 4/22/03	12:15	MD1XS9		
D1XT1	Subsurface Soil (24"-36")/ Kevin Simmons	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-01-SB	S: 4/22/03	9:15	MD1XT1	 	
D1XT2	Surface Soil (0"-6")/ Kevin Simmons	ĽG	MOICA (21), VOA-s (21)	(Ice Only) (10)	GS-01-SS	S: 4/22/03	9:00	MD1XT2		
D1XT3	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-13-SS	S: 4/22/03	13:15	MD1XT3		
D1XT4	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	G\$-13-SB	S: 4/22/03	13:25	MD1XT4		

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:		
	D1XS7, D1XT2	Word Robbill - Photon			
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?		
MOICA = CLP %Moistu	re/Calcareous (soil), VOA-s = CLP TCL Volatiles (soil)				

TR Number: 4-252433676-042203-0001 REGION CO

Case No:

DAS No:

31635

Region: Sampler Signature: Chain of Custody Record Date Shipped: 4/22/03 Project Code: 03-0474 50102D04ZZQB00 FedEx Carrier Name: Account Code: Relinquished By (Date / Time) Received By (Date / Time) Airbill: 836362682006 CERCLIS ID: Shipped to: Liberty Analytical Spill ID: Corporation 2 Site Name/State: 501 Madison Avenue Gulf States Creosoting/MS Cary NC 27513 Project Leader: Brian Striggow 3 (919) 379-4100 Screening Site Investigation Action:

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No / PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT	INORGANIC SAMPLE No.	QC Type
D1XT5	Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-04-SD	S: 4/22/03	11:35	MD1XT5	-
D1XT6	Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-05-SD	S: 4/22/03	12:15	MD1XT6	<b></b>
D1XT7	Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-07-SD	S: 4/22/03	12:45	MD1XT7	- <del>-</del>
D1XT8	Surface Soil (0"-6")/	L/G .	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-23-SS	S: 4/22/03	13:55	MD1XT8	<b></b>
D1XT9	Brian Striggow Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-15-SS	S: 4/22/03	13:40	MD1XT9	· 
D1XW0	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(ice Only) (4)	GS-15-SB	S: 4/22/03	13:50	MD1XW0	
D1XW	Surface Soil (0"-6")/ Kevin Simmons	L/G	MOICA (21), VOA-s (21)	(ice Only) (4)	GS-02-SS	S: 4/22/03	13:13	MD1XW1	<b></b>
D1XW2	Subsurface Soil (24"-36")/ Kevin Simmons	L/G	MOICA (21), VOA-s (21)	(ice Only) (4)	GS-02-SB	S: 4/22/03	. 13:34	MD1XW2	
D1XW3	Surface Soil (0"-6")/	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-14-SS	S: 4/22/03	14:40	MD1XW3	
D1XW4	Subsurface Soil (6"-12")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-14-SB	S: 4/22/03	14:50	MD1XW4	<del></del>

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	D1XS7, D1XT2	In domo Religion 2	
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?
MOICA = CLP %Moistu	ire/Calcareous (soil), VOA-s = CLP TCL Volatiles (soil)		

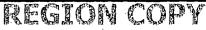
TR Number: 4-252433676-042203-0001

Sampling Co:

**USEPA SESD** 

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Case No:

DAS No:

31635

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Region: Project Code:	4		Date Shipped:	4/22/03	Chain of Custod	y Record	Sampler Signature:	
Account Code:	03-0474 50102D04ZZQB00		Carrier Name: Airbill:	FedEx 836362682006	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID: Spill ID:			Shipped to:	Liberty Analytical	1			_
Site Name/State:	Gulf States Creosoting/MS		<u>.</u>	Corporation 501 Madison Avenue	2			
Project Leader:	Brian Striggow	1		Cary NC 27513 (919) 379-4100	3			
Action: Sampling Co:	Screening Site Investigation				4		<del></del>	

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT	INORGANIC SAMPLE No.	QÇ Type
D1XW5	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-10-SS	S: 4/22/03	15:05	MD1XW5	<del></del>
D1XW6	Subsurface Soil (6"-12")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-10-SB	S: 4/22/03	15:15	MD1XW6	•
D1XW8	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-11-SS	S: 4/22/03	16:00	MD1XW8	
D1XW9	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-11-SB	S: 4/22/03	16:10	MD1XW9	
D1XX0	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-09-SS	S: 4/22/03	16.15	MD1XX0	······································
D1XX1	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-09-SB	S: 4/22/03	16:25	MD1XX1	<del></del> ·
D1XX2	Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-06-SD	S: 4/22/03	15:15	MD1XX2	<del></del>
D1XX3	Sediment/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-08-SD	S: 4/22/03	15:50	MD1XX3	<b></b>
D1XX4	Surface Soil (0"-6")/ Kevin Simmons	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-22-SS	S: 4/22/03	16:42	MD1XX4	••
D1XX5	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-06-SS	S: 4/22/03	16:40	MD1XX5	<u>-</u>

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Bampler Signature(s):	Chain of Custody Seal Number:
	D1XS7, D1XT2	In drow toldo Change	
Analysis Key:	Concentration: £ = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?
MOICA = CLP %Moistu	re/Calcareous (soil), VOA-s = CLP TCL Volatiles (soil)		

TR Number: 4-252433676-042203-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs. TR Number:

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### **Organic Traffic Report & Chain of Custody Record**

Case No:

31635

DAS No:

Region: Project Code:	4	Date Shipped:	4/22/03	Chain of Custod	y Record	Sampler Signature:	
Account Code:	03-0474 50102D04ZZQB00	Carrier Name: Airbill:	FedEx 836362682006	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID: Spill ID:		Shipped to:	Liberty Analytical	1			
Site Name/State:	Gulf States Creosoting/MS		Corporation 501 Madison Avenue	2			
Project Leader: Action:	Brian Striggow Screening Site Investigation	į	Cary NC 27513 (919) 379-4100	3			
Sampling Co:	USEPA SESD		,	4			

				1					
ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No.J PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT TE/TIME	INORGANIC SAMPLE No.	QC Type
D1XX6	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-06-SB	S: 4/22/03	16:50	MD1XX6	
D1XX7	Surface Soil (0"-6")/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-24-SS	S: 4/22/03	16:35	MD1XX7	
D1XX8	Trip Blank - Soil/ Brian Striggow	L/G	VDA-s (21)	(Ice Only) (2)	GS-01-TB	S: 4/22/03	17:54		Field QC

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	D1XS7, D1XT2	Roll 52 Offeren	
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?
MOICA = CLP %Moistu	re/Calcareous (soil), VOA-s = CLP TCL Volatiles (soil)		

TR Number: 4-252433676-042203-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs.
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F2V5.1.045 Page 5 of 5

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Case No:

DAS No:

31635

Región: Project Code:	4 03-0474		Date Shipped: Carrier Name:	4/23/03 FedEx	Chain of Custod	y Record	Sampler Signature:	
Account Code:	50102D04ZZQB00		Airbill:	836362681960	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:			Shipped to:	Liberty Analytical	1 B. Striagen	4-23-03/17:00	·	
Spill ID:		•		Corporation	D-SITIGHEN	120071100	<del> </del>	
Site Name/State:	Gulf States Creosoting/MS			501 Madison Avenue	2			. ,
Project Leader:	Brian Striggow			Cary NC 27513 (919) 379-4100	3			
Action:	Screening Site Investigation!						<del> </del>	
Sampling Co:	USEPA SESD				4			-·

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION	SAMPLE ( DATE/		INORGANIC SAMPLE No.	QC Type
D1XX9	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(ice Only) (4)	GS-12-SS	S: 4/23/03	7:40	MD1XX9	<del></del>
D1XY0	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-12-SB	S: 4/23/03	. 7:50	MD1XY0	
D1XY1	Surface Soil (0"-6")/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (10)	GS-03-SS	S: 4/23/03	7:40	MD1XY1	
D1XY2	Subsurface Soil (24"-36")/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-03-SB	S: 4/23/03	7:55	MD1XY2	<del>-</del>
D1XY3	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-08-SS	S: 4/23/03	8:1.0	MD1XY3	<del>-</del> '
D1XY4	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-08-SB	S: 4/23/03	8:20	MD1XY4	
D1XY5	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-07-SS	S: 4/23/03	8:35	MD1XY5	• <del>-</del> -,
D1XY6	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-07-SB	S: 4/23/03	8:45	MD1XY6	<u></u> ·
D1XY7.	Surface Soil (0"-6")/ Bobby Lewis	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-04-SS	S: 4/23/03	8:30	MD1XY7	
D1XY8	Subsurface Soil (24"-36")/ Bobby Lewis	L/G	MOICA (21). VOA-s (21)	(Ice Only) (4)	GS-04-SB	S: 4/23/03	8:40	MD1XY8	

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
Completer N		DR my Resse	
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?
MOICA = CLP %Moistu	re/Calcareous (soil), VOA-s = CLP TCL Volatiles (soil)		

TR Number: 4-252433676-042303-0002
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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F2V5.1.045 Page 1 of 2

Case No:

DAS No:

31635

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Region: Project Code:	4 03-0474	l l	Shipped: 4/23/03 r Name: FedEx		Chain of Custod	ly Record	Sampler Signature:	
Account Code:	50102D04ZZQB00	Airbiii:		960	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:		Shippe			1 Brsminous	4-13-03 17:00		
Spill ID:	:		Corporatio	•	DrStringer	1-19-03 17:00		
Site Name/State:	Gulf States Creosoting/MS	•	501 Madis		2		[	
Project Leader:	Brian Striggow	i	Cary NC 2 (919) 379-		3			
Action:	Screening Site Investigation	• 1	(919) 379-	100	ļ <u> </u>			<del></del>
Sampling Co:	USEPA SESD	.		•	4	•		

Camping Co.	UOLFA OL	30		1	· · · · · · · · · · · · · · · · · · ·				
ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT TE/TIME	INORGANIC SAMPLE No.	QC Type
D1XY9	Surface Soil (0"-6")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(Ice Only) (4)	GS-05-SS	S: 4/23/03	9:15	MD1XY9	
D1XZ0	Subsurface Soil (24"-36")/ Dan Thoman	L/G	MOICA (21), VOA-s (21)	(ice Only) (4)	GS-05-SB	S: 4/23/03	9:25	MD1XZ0	<del></del>
D1XZ2	Trip Blank - Soil/ Brian Striggow	L/G	VOA-s (21)	(Ice Only) (2)	GS-02-TS	S: 4/23/03	15:57		Field QC

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
		DRam Rello	
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?
MOICA = CLP %Moistu	ire/Calcareous (soil), VOA-s = CLP TCL Volatiles (soil)		

TR Number: 4-252433676-042303-0002
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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REGION COPY



Case No:

31635

DAS No:

		<u>:</u>						· · · · · · · · · · · · · · · · · · ·
Region:	4		Date Shipped:	4/25/03	Chain of Custoo	ly Record	Sampler B	1
Project Code:	03-0474		Carrier Name:	FedEx			Signature:	,
Account Code:	50102D04ZZQB00		Airbill:	836362681948	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:			Shipped to:		105./-	11 20 /		
Spill ID:			Shipped to:	Liberty Analytical Corporation	100,1	4-2503/14:40	ļ	<del></del>
Site Name/State:	Gulf States Creosoting/MS	:	<u>'</u>	501 Madison Avenue	2		1	
Project Leader:	Brian Striggow	•	l	Cary NC 27513	3			
Action:	Screening Site Investigation			(919) 379-4100	3			
Sampling Co:	USEPA SESD			•	4			

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		COLLECT E/TIME	INORGANIC SAMPLE No.	QC Type	
D1XR4	Sediment/ Bobby Lewis	. L/G	Ext/PAH (s (21)	(ice Only) (1)	GS-01-SD	S: 4/22/03	9:00	MD1XR4		:
D1XR5	Subsurface Soil (24"-36")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-18-SB	S: 4/22/03	8:50	. MD1XR5		
D1XR6	Dan Thoman Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(ice Only) (1)	GS-18-SS	S: 4/22/03	8:40	MD1XR6	· ·	
D1XR7	Dan Thoman Sübsurface Soil (24"-36")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	·G\$-19-SB	S: 4/22/03	9:35	MD1XR7		
D1XR8	Dan Thoman Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-19-SS	S: 4/22/03	9:30	MD1XR8	-2	
D1XR9	Dan Thoman Sediment/ Bobby Lewis	L/G	Ext/PAH (s (21)	(Ice Only) (1)	GS-02-SD	S: 4/22/03	9:50	MD1XR9		
D1XS0	Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-17-SS	S: 4/22/03	10:05	MD1XS0	·	
D1XS1	Dan Thoman Subsurface Soil (24"-36")/	L/G	BNA/PEST-s (21)	(ice Only) (1)	GS-17-SB	S: 4/22/03	10:20	MD1XS1	<u>:</u>	
D1XS2	Dan Thoman Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-20-SS	S: 4/22/03	10:45	MD1XS2		:
D1XS3	Dan Thoman Subsurface Soil (24"-36")/ Dan Thoman	. L∕G	BNA/PEST-s (21)	(Ice Only) (1)	GS-20-SB	S: 4/22/03	10:55	MD1XS3	<b>~~</b>	i :

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:	
	D1XT2		·	
Analysis Key:	Concentration: L = Low M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?	
BNA/PEST-s = CLP TC	Semivolatiles and Pesticides , Ext/PAH (s = CLP BNA/PEST	/LowPAH (sediment.)		

TR Number: 4-252433676-042503-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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Case No:

DAS No:

31635

Region: Project Code: Account Code:	4 03-0474 50102D04ZZQB00	Date Shipped: Carrier Name: Airbill:	4/25/03 FedEx 836362681948	Chain of Custody Re	ocord (Date / Time)	Sampler Signature: 3	(Date / Time)
CERCLIS ID: Spill ID:	:	Shipped to:	Liberty Analytical	18.6 4-29.	03 /14:40		
Site Name/State:	Gulf States Creosoting/MS		Corporation 501 Madison Avenue	2			•
Project Leader: Action:	Brian Striggow Screening Site Investigation		Cary NC 27513 (919) 379-4100	3			
Sampling Co:	USEPA SESD		•	4			

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT TE/TIME	INORGANIC SAMPLE No.	QC Type
D1XS4	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-20-SD	S: 4/22/03	10:45	MD1XS4	<del></del>
D1XS5	Sediment/ Bobby Lewis	L/G	Ext/PAH (s (21)	(Ice Only) (1)	GS-03-SD	S: 4/22/03	10:30	MD1XS5	
D1XS6	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(ice Only) (1)	GS-16-SS	S: 4/22/03	11:30	MD1XS6	
D1XS7	Subsurface Soil (24"-36")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-16-SB	S: 4/22/03	11:40	MD1XS7	<del></del>
D1XS8	Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-21-SS	S: 4/22/03	12:05	MD1XS8	
D1XS9	Dan Thoman Subsurface Soil (24"-36")/ Dan Thoman	L/G	BNA/PEST-s (21)	(ice Only) (1)	GS-21-SB	S: 4/22/03	12:15	MD1XS9	- :
D1XT1-	Subsurface Soil (24"-36")/ Kevin Simmons	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-01-SB	S: 4/22/03	9:15	MD1XT1	
D1XT2	Surface Soil (0"-6")/ Kevin Simmons	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-01-SS	S: 4/22/03	9:00	MD1XT2	
D1XT3	Surface Soll (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-13-SS	S: 4/22/03	13:15	MD1XT3	<u></u>
D1XT4	Subsurface Soil (24"-36")/ Dan Thoman	L∕G	BNA/PEST-s (21)	(ice Only) (1)	GS-13-SB	S: 4/22/03	13:25	MD1XT4	

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	D1XT2		
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?
BNA/PEST-s = CLP TC	L Semivolatiles and Pesticides , Ext/PAH (s = CLP BNA/PE	ST/LowPAH (sediment.)	

TR Number: 4-252433676-042503-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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Case No:

DAS No:

31635

	<del></del>					1		
Region: Project Code:	4		Date Shipped: Carrier Name:	4/25/03	Chain of Cust	ody Record	Sampler Signature: 03	
Account Code:	03-0474 50102D04ZZQB00		Airbili:	FedEx 836362681948	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:			Shipped to:	Liberty Analytical	13.2	4-25-03/14:40		.
Spill ID: Site Name/State:	Gulf States Creosoting/MS			Corporation 501 Madison Avenue	2	., ., ., ., ., ., ., ., ., ., ., ., ., .		
Project Leader:	Brian Striggow	į		Cary NC 27513 (919) 379-4100	3	<del> </del>		
Action:	Screening Site Investigation			(010) 010 4100	4			

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT	INORGANIC SAMPLE No.		QC Type	:
D1XT5	Sediment/ Bobby Lewis	L/G	Ext/PAH (s (21)	(Ice Only) (1)	GS-04-SD	S: 4/22/03	11:35	MD1XT5	• .		
D1XT6	Sediment/ Bobby Lewis	.L/G	Ext/PAH (s (21)	(Ice Only) (1)	GS-05-SD	S: 4/22/03	12:15	MD1XT6			i !
D1XT7	Sediment/ Bobby Lewis	L/ <sub>.</sub> G	Ext/PAH (s (21)	(Ice Only) (1)	GS-07-SD	S: 4/22/03	12:45	MD1XT7		<del></del>	
D1XT8	Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-23-SS	S: 4/22/03	13:55	MD1XT8			
D1XT9	Brian Striggow Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-15-SS	S: 4/22/03	13:40	MD1XT9			
D1XW0	Dan Thoman Subsurface Soll (24"-36")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-15-SB	S: 4/22/03	13:50	MD1XW0			; }`.
D1XW1	Surface Soil (0"-6")/ Kevin Simmons	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-02-SS	S: 4/22/03	13:13	MD1XW1			
D1XW2	Subsurface Soil (24"-36")/ Kevin Simmons	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-02-SB	S: 4/22/03	13:34	MD1XW2			İ
D1XW3	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-14-SS	S: 4/22/03	14:40	MD1XW3			
D1XW4	Subsurface Soil (6"-12")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-14-SB	S: 4/22/03	1,4:50	MD1XW4			;

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	D1XT2		
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?
BNA/PEST-s = CLP TCL	Semivolatiles and Pesticides , Ext/PAH (s = CLP BNA/PEST	/LowPAH (sediment.)	

TR Number: 4-252433676-042503-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

31635 Case No: DAS No:

Region: Project Code:	4 03-0474	Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custo	dy Record	Sampler Signature:	3,
Account Code:	50102D04ZZQB00	Airbill:	836362681948	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:		Shipped to:	Liberty Analytical	15.6	4-25-03/14:40		
Spill ID:			Corporation		1-23-05/19.40	<del> </del>	
Site Name/State:	Gulf States Creosoting/MS	•	501 Madison Avenue	2		<u> </u>	<u> </u>
Project Leader:	Brian Striggow	1	Cary NC 27513 (919) 379-4100	3			
Action:	Screening Site Investigation		(313) 673-4100			<del> </del>	<del></del>
Sampling Co:	USEPA SESD			4	·	<u> </u>	

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		COLLECT E/TIME	INORGANIC SAMPLE No.	QC Type	
D1XW5	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-10-SS	S: 4/22/03	15:05	MD1XW5		
D1XW6	Subsurface Soil (6"-12")/ Dan Thoman	L/G	BNA/PEST-s (21)	(ice Only) (1)	GS-10-SB	S: 4/22/03	15:15	MD1XW6	-	
D1XW8	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(ice Only) (1)	GS-11-SS	S: 4/22/03	16:00	MD1XW8		
D1XW9	Subsurface Soil (24"-36")/ Dan Thoman	IJG	BNA/PEST-s (21)	(Ice Only) (1)	GS-11-SB	S: 4/22/03	16:10	MD1XW9	:	
D1XX0	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(ice Only) (1)	GS-09-SS	S: 4/22/03	16:15	MD1XX0	r i d <u>u</u> a di di G	
D1XX1	Subsurface Soil (24"-36")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-09-SB	S: 4/22/03	16:25	MD1XX1	<u>-</u>	
D1XX2	Sediment/ Bobby Lewis	L/G	Ext/PAH (s (21)	(lcé Only) (1)	GS-06-SD	S: 4/22/03	15:15	MD1XX2		
D1XX3	Sedlment/ Bobby Lewis	L/G	Ext/PAH (s (21)	(Ice Only) (1)	GS-08-SD	S: 4/22/03	15:50	MD1XX3	-	
D1XX4	Surface Soil (0"-6")/ Kevin Simmons	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-22-SS	S: 4/22/03	16:42	MD1XX4		
D1XX5	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-06-SS	S: 4/22/03	16:40	MD1XX5	·	

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:		
	D1XT2				
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?	Γ.	
BNA/PEST-s = CLP TCL	Semivolatiles and Pesticides , Ext/PAH (s = CLP BNA/PEST)	LowPAH (sediment.)	•	:	

TR Number: TR Number: 4-252433676-042503-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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Case No:

31635

DAS No:

Region: Project Code:	4 03-0474	·····	Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custo	dy Record	Sampler Signature:	31
Account Code:	50102D04ZZQB00		Airbill:	836362681948	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:	•		Shipped to:	Liberty Analytical	1 18.1/	4-25.03/14-20		:
Spill ID:			.,	Corporation	10, 0	(2) - 1/ 19:00		1
Site Name/State:	Gulf States Creosoting/MS			501 Madison Avenue	2			•
Project Leader:	Brian Striggow			Cary NC 27513 (919) 379-4100	3		· ·	
Action:	Screening Site Investigation			(919) 379-4100		·	<del> </del>	<u> </u>
Sampling Co:	USEPA SESD				4			

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		COLLECT E/TIME	INORGANIC SAMPLE No.	QC Type	
D1XX6	Subsurface Soll (24"-36")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-06-SB	S: 4/22/03	16:50	MD1XX6		:
D1XX7	Surface Soil (0"-6")/ Bobby Lewis	L/G	BNA PEST-s (21)	(ice Only) (1)	GS-24-SS	S: 4/22/03	16:35	MD1XX7		
D1XX9	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-12-SS	S: 4/23/03	7:40	MD1XX9	<del>-,</del>	!
D1XY0	Subsurface Soil (24"-36")/ Dan Thoman	ĽG	BNA/PEST-s (21)	(Ice Only) (1)	GS-12-SB	S: 4/23/03	7:50	MD1XY0	 ,	
D1XY1	Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-03-SS	S: 4/23/03	7:40	MD1XY1		.] .
D1XY2	Bobby Lewis Subsurface Soil (24"-36")/	IJĠ	BNA/PEST-s (21)	(Ice Only) (1)	GS-03-SB	S: 4/23/03	7:55	MD1XY2	<del></del> .	
D1XY3	Bobby Lewis Surface Soil (0"-6")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-08-SS	S: 4/23/03	8:10	MD1XY3		
D1XY4	Dan Thoman Subsurface Soil (24"-36")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-08-SB	S: 4/23/03	8:20	MD1XY4	• 	
D1XY5	Dan Thoman Surface Soil (0"-6")/	ĽG	BNA/PEST-s (21)	(Ice Only) (1)	GS-07-SS	S: 4/23/03	8:35	MD1XY5	· 	
D1XY6	Dan Thoman Subsurface Soil (24"-36")/ Dan Thoman	ĽG	BNA/PEST-s (21)	(ice Only) (1)	GS-07-SB	S: 4/23/03	8:45	MD1XY6		

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	D1XT2		i
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High .	Type/Designate: Composite = C, Grab = G	Shipment iced?
BNA/PEST-s = CLP TCL	Semivolatiles and Pesticides , Ext/PAH (s = CLP BNA/PES	T/LowPAH (sediment.)	•

TR Number: 4-252433676-042503-0001

PR provides preliminary results. Requests for preliminary results will increase analytical costs.

Send Copy to: Sample Management Office, 2000 Edmund Halley Dr., Reston, VA. 20191-3400 Phone 703/264-9348 Fax 703/264-9222

F2V5.1.045 Page 5 of 6



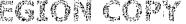
Case No: 31635 DAS No:

Region: Project Code:	4 03-0474	Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custo	dy Record	Sampler Signature:	~
Account Code:	50102D04ZZQB00	Alrbill:	836362681948	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:	**	Shipped to:	Liberty Analytical	18.0	4-25-03/14:40		
Spill ID:	•	} ''	Corporation		12707/17	<del> </del>	<del></del>
Site Name/State:	Gulf States Creosoting/MS	1	501 Madison Avenue	2		·	<u> </u>
Project Leader:	Brian Striggow	·	Cary NC 27513 (919) 379-4100	3			
Action:	Screening Site Investigation	J	(313) 313-4100			ļ	
Sampling Co:	USEPA SESD		·	4			i

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT TE/TIME	INORGANIC SAMPLE No.	QC Type	
D1XY7	Surface Soil (0"-6")/ Bobby Lewis	IJĠ	BNA/PEST-s (21)	(Ice Only) (1)	GS-04-SS	S: 4/23/03	8:30	MD1XY7		: .
D1XY8	Subsurface Soil (24"-36")/ Bobby Lewis	r\è	BNA/PEST-s (21)	(Ice Only) (1)	GS-04-SB	S: 4/23/03	8:40	MD1XY8	, <del></del> .	
D1XY9	Surface Soil (0"-6")/ Dan Thoman	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-05-SS	S: 4/23/03	9:15	MD1XY9		!
D1XZ0	Subsurface Soil (24"-36")/	L/G	BNA/PEST-s (21)	(Ice Only) (1)	GS-05-SB	S: 4/23/03	9:25	MD1XZ0		. !

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC: D1XT2	Additional Sampler Signature(s):	Chain of Custody Seal Number:	
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment iced?	
BNA/PEST-s = CLP TCL	Semivolatiles and Pesticides , Ext/PAH (s = CLP BNA/PEST/	LowPAH (sediment.)		

TR Number: 4-252433676-042503-0001
PR provides preliminary results. Requests for preliminary results will increase analytical costs. Send Copy to: Sample Management Office, 2000 Edmund Halley Dr., Reston, VA. 20191-3400 Phone 703/264-9348 Fax 703/264-9222



	Express USA Airbill Fraction of Company A 3 L 3	362687448
1	1 From Playes program And Proses had Date 4/25/C-3 Sender's FedEx Account Number 2/48-6	184.5
	Sender's Brian Striggin Phone	706 355-8619
	COMPANY USEPA SESD	
	Address 980 College Star	tin Al
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2	2 Your Internal Billing Reference (1873) that the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state of the state o	L .
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FedEx 2Day Second business day FedEx Envelope rate not available.	FedEx Express Saver Third business day Minimum charge: One-pound (als	_
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Case No:

DAS No:

31635

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Reģion: Project Code:	4		Date Shipped:	4/25/03	Chain of Custody	/ Record	Sampler Signature:	e
Account Code:	03-0474 50102D04ZZQB00	1	Carrier Name:	FedEx	Relinguished By	(Date / Time)	Received By	(Date / Time)
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Split ID:		:	l ''	1544 Sawdust Road			<del>                                     </del>	
Site Name/State:	Gulf States Creosoting/MS	I		Suite 505	2	_		
Project Leader:	Brian Striggow	:	1	The Woodlands TX 77380	3			
Action:	Screening Site Investigation		i	(281) 292-5277	<u> </u>		<u> </u>	
			l .		1			
Sampling Co:	USEPA SESD		1 .		[ 4	•		

ORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT TE/TIME	INORGANIC SAMPLE No.	QC Type
D1XT0	Ground Water/ Kevin Simmons	L/G	BNA/PEST-w (21), VOA-w (21)	(Ice Only) (14)	GS-01-GW	S: 4/22/03	11:05	MD1XT0	
D1XW7	Ground Water/ Kevin Slmmons	L/G	BNA/PEST-w (21), VOA-w (21)	(Ice Only) (6)	GS-02-GW	S: 4/22/03	14:55	MD1XW7	. <del></del>
D1XZ1	Ground Water/ Brian Striggow	L/G	BNA/PEST-w (21), VOA-w (21)	(ice Only) (6)	GS-03-GW	S: 4/23/03	9:55	MD1XZ1	. <del></del>
D1XZ3	Ground Water/ Brian Striggow	L/G	BNA/PEST-w (21), VOA-w (21)	(Ice Only) (6)	GS-04-GW	S: 4/23/03	13:20	MD1XZ3	<del></del>
D1XZ4	Ground Water/ Brian Striggow	. L/G	BNA/PEST-w (21), VOA-w (21)	(ice Only) (6)	GS-05-GW	S: 4/23/03	11:45	MD1XZ4	<del></del> .

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:	
!	D1XT0		·	
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?	
BNA/PEST-w = CLP TC	L Semivolatiles and Pesticides , VOA-w = CLP TCL Volatiles (v	vater)		

TR Number: 4-252433676-042503-0002
PR provides preliminary results. Requests for preliminary results will increase analytical costs. TR Number:





Case No:

DAS No:

31635

Region: Project Code:	4 03-0474		Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custod	y Record	Sampler Signature:	
Account Code:	50102D04ZZQB00	1	Airbill:	836362681981	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:			Shipped to:	Sentinel Inc.	1 3. 12	4-25-03 /14.40		
Spill ID:				116 Washington Street,	0.72	עבידות כשינגי		
Site Name/State:	Guif States Creosoting/MS			NE.	<b>1</b> 2		·	
Project Leader:	Brian Striggow			Huntsville AL 35801 (256) 534-9800	3			
Action:	Screening Site Investigation			(230) 334-9600		<del></del>	·	
Sampling Co:	USEPA SESD		,		4			

INORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT	ORGANIC SAMPLE No.	QC Typ <del>d</del>
MD1XR3	Ground Water/ Brian Striggow	L/G	TM-w (21)	(HNO3) (1)	GS-06-GW	S: 4/24/03	.13:45	·	
MD1XR4	Sediment/ Bobby Lewis	ĽG.	TM-s (21)	(Ice Only) (1)	GS-01-SD	S: 4/22/03	9:00	D1XR4	
MD1XR5	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(ice Only) (1)	GS-18-SB	S: 4/22/03	8:50	D1XR5	· <del>_</del> .
MD1XR6	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-18-SS	S: 4/22/03	8:40	D1XR6	
MD1XR7	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-19-SB	S: 4/22/03	9:35	D1XR7	 -
MD1XR8	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-19-SS	S: 4/22/03	9:30	D1XR8	<del>-</del>
MD1XR9	Sediment/ Bobby Lewis	L/G	TM-s (21)	(Ice Only) (1)	GS-02-SD	S: 4/22/03	9:50	D <sub>1</sub> 1XR9	<del>-</del> ·
MD1XS0	Surface Soil (0"-6")/ Dan Thoman	ĽG .	TM-s (21)	(ice Only) (1)	GS-17-SS	S: 4/22/03	10:05	D1XS0	, <del></del>
MD1XS1	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-17-SB	S: 4/22/03	10:20	D1XS1	<del></del>
MD1XS2	Surface Soil (0"-6")/	L/G	TM-s (21)	(Ice Only) (1)	GS-20-SS	S: 4/22/03	10:45	D1XS2	<u>.</u>

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:	
	MD1XT0, MD1XT2			
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?	
TM-s = CLP TAL Total N	letals (soil), TM-w = CLP TAL Total Metals (water)			

TR Number: 4-252433676-042503-0003
PR provides preliminary results. Requests for preliminary results will increase analytical costs.



## USEPA CONTRACT Laboratory Program Inorganic Traffic Report & Chain of Custody Record

Case No:

31635

DAS No:

Region: Project Code:	4 03-0474	Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custod	y Record	Sampler Signature:	e-
Account Code:	50102D04ZZQB00	Airbill:	836362681981	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:		Shipped to:	Sentinel Inc.	18.2	4.25.03/14:40		
Spill ID:	i		116 Washington Street,	100	12703714.40		<del></del>
Site Name/State:	Gulf States Creosoting/MS		NE	2		١	
Project Leader: Action:	Brian Striggow Screening Site Investigation		Huntsville AL 35801 (256) 534-9800	3.			
Sampling Co:	USEPA SESD			4			

INORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		COLLECT E/TIME	ORGANIC SAMPLE No.	QC Type
MD1XS3	Subsurface Soil (24"-36")/ Dan Thoman	ĽG	TM-s (21)	(Ice Only) (1)	GS-20-SB	S: 4/22/03	10:55	D1XS3	-
MD1XS4	Surface Soil (0"-6")/ Dan Thoman	L∕G	TM-s (21)	(Ice Only) (1)	GS-20-SD	S: 4/22/03	10:45	D1XS4	<u></u>
MD1XS5	Sediment/ Bobby Lewis	L/G	TM-s (21)	(Ice Only) (1)	GS-03-SD	S: 4/22/03	10:30	D1XS5	 ·
MD1XS6	Surface Soil (0"-6")/ Dan Thoman	L∕G	TM-s (21)	(Ice Only) (1)	GS-16-SS	S: 4/22/03	11:30	D1XS6	••• •
MD1X\$7	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-16-SB	S: 4/22/03	11:40	D1XS7	<del></del>
MD1XS8	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(ice Only) (1)	GS-21-SS	S: 4/22/03	12:05	D1XS8	·
MD1XS9	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-21-SB	S: 4/22/03	12:15	D1XS9	. <del></del>
MD1XTO	Ground Water/ Kevin Simmons	<b>L/G</b>	TM-w (21)	(HNO3) (2)	GS-01-GW	S: 4/22/03	11:05	D1XT0	
MD1XT1	Subsurface Soil (24"-36")/ Kevin Simmons	L/G	TM-s (21)	(Ice Only) (1)	GS-01-SB	S: 4/22/03	9:15	D1XT1	<b>-</b> •
MD1XT2	Surface Soil (0"-6")/ Kevin Simmons	L/G	TM-s (21)	(Ice Only) (1)	GS-01-SS	S: 4/22/03	9:00	D1XT2	-

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	MD1XT0, MD1XT2		
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?
TM-s = CLP TAL Total M	etals (soil), TM-w = CLP TAL Total Metals (water)		

TR Number: 4-252433676-042503-0003
PR provides preliminary results. Requests for preliminary results will increase analytical costs. TR Number:



Case No:

DAS No:

31635

1635

Region: Project Code:	4 03-0474		Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custod	y Record	Sampler Signature: B	~
Account Code:	50102D04ZZQB00	_	Airbill:		Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:	-		Shipped to:	Sentinel Inc.	1 B. 1	4.25-03/14.40		
Spill ID:	4	•	••	116 Washington Street,		64700/11.75		
Site Name/State:	Gulf States Creosoting/MS			NE	2			·
Project Leader:	Brian Striggow			Huntsville AL 35801 (256) 534-9800	3			
Action:	Screening Site Investigation			(250) 554-9600			<u> </u>	
Sampling Co:	USEPA SESD	·		·	4		ŀ	

INORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E COLLECT FE/TIME	ORGANIC SAMPLE No.	QC Type
MD1XT3	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(ice Only) (1)	GS-13-SS	S: 4/22/03	13:15	D1XT3	_
MD1XT4	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-13-SB	S: 4/22/03	13:25	D1XT4	- •
MD1XT5	Sediment/ Bobby Lewis	L/G	TM-s (21)	(ice Only) (1)	GS-04-SD	S: 4/22/03	11:35	D1XT5	<del></del>
MD1XT6	Sediment/ Bobby Lewis	L/G	TM-s (21)	(Ice Only) (1)	GS-05-SD	S: 4/22/03	12.15	D1XT6	
MD1XT7	Sediment/ Bobby Lewis	L/G	TM-s (21)	(Ice Only) (1)	GS-07-SD	S: 4/22/03	12:45	D1XT7	<b></b>
MD1XT8	Surface Soil (0"-6")/ Brian Striggow	. L/G	TM-s (21)	(Ice Only) (1)	GS-23-SS	S: 4/22/03	13:55	D1XT8	
MD1XT9	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-15-SS	S: 4/22/03	13:40	D1XT9	. <del></del>
MD1XW0	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-15-SB	S: 4/22/03	13:50	D1XW0	<u>.</u>
MD1XW1	Surface Soil (0"-6")/ Kevin Simmons	L/G	TM-s (21)	(ice Only) (1)	GS-02-SS	s. 4/22/03	13:13	D1XW1	<b></b>
MD1XW2	Subsurface Soil (24"-36")/ Kevin Simmons	ĽG	TM-s (21)	(Ice Only) (1)	` GS-02-SB	S: 4/22/03	13:34	D1XW2	

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:		
	MD1XT0, MD1XT2		•		
Analysis Key:	Concentration: L = Low,M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?		
TM-s = CLP TAL Total N	etals (soil), TM-w = CLP TAL Total Metals (water)				

TR Number: 4-252433676-042503-0003
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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## USERA CONTRACT Laboratory Program Inorganic Traffic Report & Chain of Custody Record

Case No:

DAS No:

31635

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		<del></del>						
Region: Project Code:	4		Date Shipped:	4/25/03	Chain of Custody	Record	Sampler Signature:	11_
· ·	03-0474	1	Carrier Name:	FedEx	<u> </u>		ļ	
Account Code:	50102D04ZZQB00	:	Airbill:	836362681981	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:		•	Shipped to:	Sentinel Inc.	181=	V 25 02/11011		
Spill ID:		:	Simpled to.	116 Washington Street.	1 p. 1	4.25.03/1440	<u> </u>	
Site Name/State:	Gulf States Creosoting/MS			NE	2			
Project Leader:	Brian Striggow	}	1	Huntsville AL 35801	3			
Action:	Screening Site Investigation	<u>:</u>		(256) 534-9800	3		<del></del>	
Sampling Co:	USEPA SESD	!	1		4	.,	}	

INORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION		E.COLLECT FE/TIME	ORGANIC SAMPLE No.	QC Type	•
MD1XW3	Surface Soil (0"-6")/ Dan Thoman	IJĠ	TM-s (21)	(Ice Only) (1)	GS-14-SS	S: 4/22/03	14:40	D1XW3		
MD1XW4	Subsurface Soil (6"-12")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-14-SB	S: 4/22/03	14:50	D1XW4		
MD1XVV5	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-10-SS	S: 4/22/03	15:05	D1XW5		
MD1XW6	Subsurface Soil (6"-12")/ Dan Thoman	L/G	TM-s (21)	(ice Only) (1)	GS-10-SB	S: 4/22/03	15:15	D1XW6		
MD1XVV7	Ground Water/ Kevin Simmons	L/G	TM-w (21)	(HNO3) (1)	GS-02-GW	S. 4/22/03	14:55	D1XW7	 ·	
MD1XW8	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-11-SS	S: 4/22/03	16:00	D1XW8		
MD1XW9	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-11-SB	S: 4/22/03	16:10	D1XW9	·	
MD1XX0	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-09-SS	S: 4/22/03	16:15	D1XX0	·	÷
MD1XX1	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-09-SB	S: 4/22/03	16:25	D1XX1	 ·	
MD1XX2	Sediment/ Bobby Lewis	ĽG	TM-s (21)	(ice Only) (1)	GS-06-SD	S: 4/22/03	15:15	.D1XX2	·	

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:		
	MD1XT0, MD1XT2	-			
Analysis Key:	Concentration: L = Low, M = Low/Medium, H ≈ High	Type/Designate: Composite = C, Grab = G	Shipment Iced?		
TM-s = CLP TAL Total N	letals (soil), TM-w = CLP TAL Total Metals (water)				

TR Number: 4-252433676-042503-0003
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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Case No:

DAS No:

31635

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Region: Project Code:	4	Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custody	Record	Sampler Signature:	
Account Code:	03-0474 50102D04ZZQB00	Alrbili:	836362681981	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:	·	Shipped to:	Sentinel Inc.	1B.W.4-	25.03 /14:40		
Spilli ID:	· !		116 Washington Street		22.00 / 17.70	<del> </del>	<del></del>
Site Name/State:	Gulf States Creosoting/MS	1	NE .	2	·		
Project Leader:	Brian Striggow		Huntsville AL 35801 (256) 534-9800	3			
Action:	Screening Site Investigation		(200) 00 1 0000			ļ	
Sampling Co:	USEPA SESD			4			

INORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION	SAMPLE COLLECT	ORGANIC SAMPLE No.	QC Type
MD1XX3	Sediment/ Bobby Lewis	L/G	TM-s (21)	(Ice Only) (1)	G\$-08-SD	S: 4/22/03 15:50	D1XX3	
MD1XX4	Surface Soil (0"-6")/ Kevin Simmons	L/G	TM-s (21)	(ice Only) (1)	GS-22-SS	S: 4/22/03 16:42	D1XX4	. <del></del>
MD1XX5	Surface Soil (0"-6")/	L/G	TM-s (21)	(Ice Only) (1)	GS-06-SS	S: 4/22/03 16:40	D1XX5	
MD1XX6	Dan Thoman Subsurface Soil (24"-36")/	L/G	TM-s (21)	(Ice Only) (1)	G\$-06-SB	S: 4/22/03 16:50	D1XX6	
MD1XX7	Dan Thoman Surface Soil (0"-6")/	IJG	TM-s (21)	(Ice Only) (1)	GS-24-SS	S: 4/22/03 16:35	D1XX7	
MD1XX9	Bobby Lewis Surface Soil (0"-6")/	L/G	TM-s (21)	(Ice Only) (1)	GS-12-SS	S: 4/23/03 7:40	D1XX9	
MD1XY0	Dan Thoman Subsurface Soil (24"-36")/	L/G	TM-s (21)	(Ice Only) (1)	GS-12-SB	S: 4/23/03 7:50	D1XY0	
MD1XY1	Dan Thoman Surface Soil (0"-6")/	L/G	TM-s (21)	(Ice Only) (1)	GS-03-SS	S: 4/23/03 7:40	D1XY1	<del></del>
MD1XY2	Bobby Lewis Subsurface Soil (24"-36")/	L/G	TM-s (21)	(Ice Only) (1)	GS-03-SB	S: 4/23/03 7:55	D1XY2	<u> </u>
MD1XY3	Bobby Lewis Surface Soil (0"-6")/	ĽG	TM-s (21)	(Ice Only) (1)	GS-08-SS	S: 4/23/03 8:10	D1XY3	
7.4	Dan Thoman				•	•		

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:	
	MD1XT0, MD1XT2			
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?	
TM-s = CLP TAL Total M	letals (soil), TM-w = CLP TAL Total Metals (water)			

TR Number: 4-252433676-042503-0003





Case No:

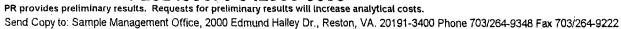
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	Thorganic Traine Repor		ustouy kecoru		DAS No:		1 \
Region: Project Code:	4 03-0474	Date Shipped: Carrier Name:	4/25/03 FedEx	Chain of Custody R	ecord	Sampler Signature:	2.1
Account Code:	50102D04ZZQB00	Airbill:	836362681981	Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID:		Shipped to:	Sentinel Inc.	1 8 1/ 4.	25.03/14.40		
Spili ID:		1	116 Washington Street,	1	3 0 3 / 14.40	<del> </del>	
Site Name/State:	Gulf States Creosoting/MS	j	NE	2			
Project Leader:	Brian Striggow	j	Huntsville AL 35801 (256) 534-9800	3			
Action:	Screening Site Investigation	1	(230) 534-9600		<del></del>	<u> </u>	
Sampling Co:	USEPA SESD	]	•	4			

INORGANIC SAMPLE No.	MATRIX/ SAMPLER	CONC/ TYPE	ANALYSIS/ TURNAROUND	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION	SAMPLE CO DATE/T		ORGANIC SAMPLE No.	QC Type	
MD1XY4	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-08-SB	S: 4/23/03	3:20	D1XY4	<del></del>	
MD1XY5	Surface Soil (0"-6")/ Dan Thoman	ĽG	TM-s (21)	(ice Only) (1)	GS-07-SS	S: 4/23/03	B:35	D1XY5	*	
MD1XY6	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-07-SB	S: 4/23/03	8:45	D1XY6		
MD1XY7	Surface Soil (0"-6")/ Bobby Lewis	L/G	TM-s (21)	(Ice Only) (1)	GS-04-SS	S: 4/23/03	8:30	D1XY7		٠
MD1XY8	Subsurface Soil (24"-36")/ Bobby Lewis	L/G	TM-s (21)	(Ice Only) (1)	GS-04-SB	S: 4/23/03	8:40	D1XY8		
MD1XY9	Surface Soil (0"-6")/ Dan Thoman	L/G	TM-s (21)	(ice Only) (1)	GS-05-SS	S: 4/23/03	9:15	D1XY9		
MD1XZ0	Subsurface Soil (24"-36")/ Dan Thoman	L/G	TM-s (21)	(Ice Only) (1)	GS-05-SB	S: 4/23/03	9:25	D1XZ0		
MD1XZ1	Ground Water/ Brian Striggow	L/G	TM-w (21)	(HNO3) (1)	GS-03-GW	S: 4/23/03	9:55	D1XZ1	· · · · · · · · · · · · · · · · · · ·	
MD1XZ3	Ground Water/ Brian Striggow	L/G	TM-w (21)	(HNO3) (1)	GS-04-GW	S: 4/23/03 1	3:20	D1XZ3	<del></del>	
MD1XZ4	Ground Water/ Brian Striggow	L/G	TM-w (21)	(HNO3) (1)	GS-05-GW	S: 4/23/03 1	1:45	D1XZ4		

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	MD1XT0, MD1XT2		
Analysis Key:	Concentration: L = Low, M = Low/Medium, H ≈ High	Type/Designate: Composite = C, Grab = G	Shipment Iced?
TM-s = CLP TAL Total N	letals (soil), TM-w = CLP TAL Total Metals (water)		

TR Number: 4-252433676-042503-0003





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Kevin Simmons

#### **USEPA Contract Laboratory Program Inorganic Traffic Report & Chain of Custody Record**

Case No:

DAS No:

31635

Region: Project Code:	4 03-0474	Date Shipped: 4/25/03  Carrier Name: FedEx		Chain of Custo	ody Record	Sampler Signature: B	
Account Code:	50102D04ZZQB00	Airbill: 836362681981		Relinquished By	(Date / Time)	Received By	(Date / Time)
CERCLIS ID: Spill ID:		Shipped to: Sentinel Inc.		1 B.N 4-25:03/14:40			
Site Name/State:	Gulf States Creosoting/MS	116 Washington NE	116 Washington Street, NE	2			
Project Leader: Action:	Brian Striggow Screening Site Investigation	Huntsville AL 356 (256) 534-9800	301	3			
Sampling Co:	USEPA SESD			4			
INORGANIC SAMPLE No.	MATRIX/ CONC/ ANALYSIS/ SAMPLER TYPE TURNAROUN	TAG No./ PRESERVATIVE/ Bottles	STATION LOCATION			RGANIC MPLE No.	QC Type

INORGANIC	MATRIX/	CONC/	ANALYSIS/	TAG No./	STATION	SAMPLE COLLECT	ORGANIC	QC	
SAMPLE No.	SAMPLER	TYPE	TURNAROUND	PRESERVATIVE/ Bottles	LOCATION	DATE/TIME	SAMPLE No.	Type	
MD1XZ5	Preservative Blank/	L/G	TM-w (21)	(HNO3) (1)	GS-01-PB	S: 4/24/03 14:00		Field QC	

Shipment for Case Complete? N	Sample(s) to be used for laboratory QC:	Additional Sampler Signature(s):	Chain of Custody Seal Number:
	MD1XT0, MD1XT2		
Analysis Key:	Concentration: L = Low, M = Low/Medium, H = High	Type/Designate: Composite = C, Grab = G	Shipment Iced?
TM-s = CLP TAL Total M	letals (soil), TM-w = CLP TAL Total Metals (water)		

TR Number: 4-252433676-042503-0003
PR provides preliminary results. Requests for preliminary results will increase analytical costs.

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Track Shipments

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Ouick Help

836362681992

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Tracking Number 836362681992 Reference Number Ship Date Delivered To Delivery Location SPRING TX Delivery Date/Time

04/26/2003 09:45 Signed For By A.4 SCINTIFIC Service Type Priority Overnight Master Tracking Number 836362681992 Total Pieces 1 of 2

04/25/2003

Shioments Track More Shipments

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Fedex Signature Proof of Delivery Get it in writing! click here.

Scan Activity Date/Time Delivered/SPRING TX 04/26/2003 09:45 Delivered/SPRING TX 04/26/2003 09:45 On FedEx vehicle for delivery/CONROE TX 04/26/2003 08:58 On FedEx vehicle for delivery/CONROE TX 04/26/2003

08:04 04/26/2003 Arrived at FedEx Destination Location/CONROE

08:03 Arrived at FedEx Ramp/HOUSTON TX 04/26/2003 04:47 Left FedEx Ramp/HOUSTON TX 04/26/2003

04:02 Left FedEx Sort Facility/MEMPHIS TN 04/26/2003 03:47 Left FedEx Ramp/ATLANTA GA 04/26/2003

00:59 Arrived at Sort Facility/MEMPHIS TN 04/25/2003 23:39 Arrived at FedEx Ramp/ATLANTA GA 04/25/2003

21:18 Left FedEx Origin Location/ATHENS GA 04/25/2003 19:23

Comments

No signature required - release waiver on

No signature required - release waiver on

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Enter your email (optional), up to three email addresses as recipients, add your message, and click on Send Email.

From To To

Add a message to this email.

Fedex USA Airbill		
Express USA AUTOUL Freder 836362681981	60m/. 0200	
1 From Process processed presentant,	4a Express Package Service	Packages up to 150 lbs. Delivery commitment may be later in some areas.
Date 1/25/03 Sender's FedEx 2148-8184-5	FedEx Priority Overnight FedEx Standard Overnigh Next business afternoon	
Sender's Brian Striggon Phone (706) 340-6352	FedEx 2Day FedEx Express Saver	
Company USEPA SESD	4b Express Freight Service	Packages over 150 lbs. Delivery commement may be later in some areas
Address 980 College Station Rd	FedEx 1Day Freight* FedEx 2Day Freight Next business day  * Califor Confirmation:	FedEx 3D ay Freight Third business day
Dept.Roor/Sure/Room	5 Packaging	• Daclared value limit \$500
city Athens State Clit ZIP 30605	FedEx Envelope*  FedEx Pak*  Michales FadEx Small Pak, FedEx	Other
2 Your Internal Billing Reference 6.9 Patroda.	Large Pak, and FedEx Sturoy Pak  6 Special Handling Inches FedEx with	!
3 To Recipient's Sample Custadian Phone 256,534-9800	SATURDAY Delivery Available ONLY for Feder Modery to Seed Freder Owngon and Feder Moyer to seed 210 cooss	ess in Section 3.  HOLD Saturday  at FedEx Location  Available ONLY for fedEx Printy Overnight and
Continue / To	Does this shipment contain dangerous goods?  One lies must be checked.	· Fedex 2Day to same trocaudns
Company SEAVINET LAC	> 24 Subbel 2 Deciaration > 40 Leading	/ry lce ry lce 9 UN 1845 \ \ \ \ \ \ \ \ \ \ \
Address 1/6 Washington Street NE	Dangerous Goods (including Dry Ice) cannot be shipped in FedEx packaging.	Cargo Aircraft Only
To "HOLD" at FedEx location, print FedEx eddress.  We cannot deliver to P.O. baxes or P.O. ZIP codes.  Address	7 Payment Bill to:  Sender Act No. n Section Recipient Third Party Act No. 19 Section I will be balled.	Credit Card Cash/Check
Papt/Roor/SulmyRoom	FedEx Acct No Credit Card No.	Exp. Date
city Hundsville State Al ZIP 35801	Total Packages Total Weight Total Decla	
	s	.00
Tryonline anipping at federacons	<sup>†</sup> Our hability is timited to \$100 unless you declare a higher value. See ba	ck for details. FedEx Use Only
By using this Airbill you agree to the service conditions on the back of this Airbill and in our current Service Guide, including terms that limit our liability.	8 Release Signature Sign to authorize delivery without obtaining signature	
Questions? Visit our Web site at fedex.com	By signing you authorize us to deliver this shipment without obtaining a s and agree to indemnify and hold us harmless from any resulting claims.	ignature 446

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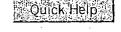
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#### Track Shipments

#### **Detailed Results**

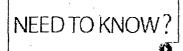
View detailed results for:

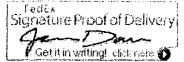


836362681981

#### Related Links

- Signature Proof
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- ▶ FedEx Wireless Solutions
- ▶ EedEx Address Checker





Tracking Number 836362681981 Reference Number

Ship Date Delivered To **Delivery Location** Delivery Date/Time 04/26/2003 10:07

Signed For By Master Tracking Number 836362681981 Total Pieces 1 of 2

04/25/2003 Ship'g/Receiv'g HUNTSVILLE AL N.MAGHSOUDLOU Service Type Priority Overnight

#### **Tracking Options**

- ★ Obtain a Signature Proof of Delivery
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  - one or more recipients Return to Multiple Piece
  - Shipments Return to Summary Results
  - Track More Shipments

Scan Activity	Date/Time	Comments
Delivered/HUNTSVILLE AL	04/26/2003 10:07	
On FedEx vehicle for delivery/MADISON AL	04/26/2003 08:20	
Arrived at FedEx Destination Location/MADISON AL	04/26/2003 08:13	
Left FedEx Sort Facility/MEMPHIS TN	04/26/2003 03:26	
Left FedEx Ramp/ATLANTA GA	04/26/2003 00:59	
Arrived at Sort Facility/MEMPHIS TN	04/25/2003 23:39	•
Arrived at FedEx Ramp/ATLANTA GA	04/25/2003 21:18	
Left FedEx Origin Location/ATHENS GA	04/25/2003 19:23	
Picked up by FedEx/ATHENS GA	04/25/2003 16:18	
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#### **UNITED STATES ENVIRONMENTAL PROTECTION AGENCY**

#### Region 4

Science and Ecosystem Support Division 980 College Station Road Athens, Georgia 30605-2720

#### **MEMORANDUM**

Date: 06/20/2003

Subject: Results of VOLATILES Sample Analysis

03-0474

**Gulf States Creosoting** 

Flowood, MS

From: Goddard, Denise

To: Striggow, Brian

Thru: QA Office

Attached are the results of analysis of samples collected as part of the subject project. If you have any questions, please contact me.

Volatiles         J         continue           6754, 6755, 6756, 6757, 6758, 6759, 6760, 6761, 6762, 6763, 6764, 6765, 6766, 6767, 6768, 6766, 6771, 6772, 6773, 6774, 6775         bromomethane         J         continue           6754, 6755, 6756, 6757, 6758, 6756, 6760, 6764, 6767, 6768, 6769, 6771, 6772, 6773, 6774, 6775         acetone, 2-butanone, 4-methyl-2-pentanone, 2-hexanone, 1,2-dibromo-3-chloropropane         J         continue           6754, 6756, 6758, 6759, 6760, 6769, 6760, 6760, 6760, 6762, 6766, 6768, 6773, 6775         acetone         J         contamblank           6754, 6756, 6758, 6758, 6758, 6764         methyl acetate         J         contamblank           6754, 6765         carbon disulfide         J         quan           6772, 6776, 6777, 6778, 6779, 6780, 6780, 6782, 6783, 6784, 6785, 6786, 6788, 6789, 6790, 6791, 6792, 6793, 6794, 6795         contamblank         blank           6772, 6776, 6777, 6798, 6799, 6794, 6795         bromomethane         J         continue           6772, 6776, 6777, 6798, 6799, 6790, 6791, 6792, 6793, 6794, 6795         bromomethane         J         continue           6772, 6776, 6777, 6778, 6779, 6780, 6780, 6781, 6782, 6783, 6786, 6782, 6783, 6786, 6782, 6783, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786, 6786,	Case Number:	31635	Project Number 03-0474	SOW	Number	OLC03.2, OLM04.2	
Affected Samples         Compound or Fraction         Used         Reference           Volatiles         Volatiles         Image: Compound or Fraction         >Site ID.</td> <td>Gulf Stat</td> <td>es Creosoting, Flowood, MS</td> <td></td> <td colspan="3">Date 6/0</td>	Site ID.	Gulf Stat	es Creosoting, Flowood, MS		Date 6/0		
6754, 6755, 6756, 6757, 6758, 6759, 6760, 6761, 6762, 6766, 6767, 6768, 6759, 6771, 6772, 6773, 6774, 6775  6754, 6755, 6756, 6757, 6758, 6759, 6760, 6764, 6767, 6768, 6769, 6771, 6772, 6773, 6774, 6775  6754, 6756, 6758, 6759, 6760, 6762, 6760, 6760, 6762, 6768, 6769, 6773, 6775  6754, 6765, 6766, 6762, 6768, 6773, 6775  6754, 6765  6754, 6765  6757, 6776, 6777, 6778, 6779, 6780, 6781, 6782, 6783, 6794, 6795  6772, 6776, 6777, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6778, 6779, 6780, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6790, 6791, 6792, 6793, 6794, 6795	Affected Samples	Flag				Reason	
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6769, 6771, 6772, 6773, 6774, 6775  6754, 6755, 6756, acetone, 2-butanone, 4-methyl-2-pentanone, J continu (6757, 6758, 6759, 2-hexanone, 1,2-dibromo-3-chloropropane)  6760, 6764, 6767, 6768, 6769, 6771, 6772, 6773, 6774, 6775  6754, 6756, 6758, acetone  6759, 6760, 6762, 6766, 6768, 6773, 6775  6754, 6765  6754, 6765  6754, 6765  6754, 6765  6754, 6765  6757, 6766, 6777, 6778, 6779, 6780, 6781, 6782, 6783, 6789, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6797, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6780, 6797, 6780, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6790, 6791, 6792, 6793, 6794, 6795  6772, 6776, 6777, 6780, 6780, 6781, 6782, 6783, 6789, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6780, 6781, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 6782, 6783, 678	6757, 6758, 6759, 6760, 6761, 6762, 6763, 6764, 6765,	bromomo	ethane	. J	continuin %D outli	g calibration er	
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· ·				
Case Number:	31635 Project Number 03-0474	SOW 1	Number	OLC03.2, OLM04.2
Site ID.	Gulf States Creosoting, Flowood, MS		Date	6/07/03
Affected Samples	Compound or Fraction	Flag <u>Used</u>	Reas	on
6770, 6787, 6811, 6812, 6813	2-butanone	J.	continuing outlier	g calibration
6788, 6789, 6790, 6791, 6792, 6793, 6794, 6795	2-butanone, 4-methyl-2-pentanone, 2-hexanone	J	continuing %D outlie	g calibration er
6779	methyl acetate	J	< quantita	tion limit
6786	benzene	J	< quantita	tion limit
6790	methyl tert-butyl ether, 1,2-dibromo-3-chloropropane	J	continuing	g calibration er
6796, 6797, 6798, 6799, 6800, 6801, 6802, 6804, 6805, 6806, 6807, 6808, 6809, 6810, 6814	acetone, 2-butanone, 4-methyl-2-pentanone, 2-hexanone	j	continuin %D outlie	g calibration er
6799, 6800, 6801, 6802, 6804, 6805, 6806, 6807, 6808, 6809, 6810, 6814	bromomethane, methyl tert-butyl ether, 1,2-dibromo-3-chloropropane	J	continuin %D outlie	g calibration er

C - Newstern	21.625	Duration AND The same	02.0474	COMA		OLC03.2,		
Case Number:	31635 Project Number 03-0474			SOW I	Number	OLM04.2		
Site ID.	Gulf State	es Creosoting, Flowo	od, MS		Date	6/07/03		
Affected Samples	Compound or Fraction				Flag <u>Used</u> <u>Reason</u>			
Extractables		·				,		
6754, 6755, 6756, 6757, 6758, 6759, 6760, 6761, 6762, 6763, 6764, 6765, 6766, 6767, 6768, 6769, 6771, 6773, 6774, 6775	butylbenz	zylphthalate		J	continuing %D outlie	calibration r		
6754, 6755, 6756, 6757, 6758	bis(2-eth	ylhexyl)phthalate		<b>J</b>	continuing %D outlie	calibration		
6759, 6760, 6761, 6762, 6763, 6764, 6765, 6766, 6767, 6768, 6769, 6771, 6773, 6774, 6775	3,3'-dich	lorobenzidine		J	continuing %D outlie	g calibration r		
6760, 6761, 6762, 6764, 6765, 6766, 6773, 6774	benzo(b)	fluoranthene, benzo(l	c)fluoranthene	J	isomers corresolved b	ould not be by GC		
6754	benzalde	hyde, benzo(g,h,i)per	ylene	J	< quantita	tion limit		
6763	benzo(g,	h,i)perylene		J	< quantita	tion limit		
6760	- carbazol	hylene, phenanthrene e, dibenzo(a,h)anthrad h,i)perylene		J	< quantita	tion limit		
6761	benzo(b)	fluoranthene, benzo(l	k)fluoranthene	J ·	< quantita	tion limit		
6762	benzo(a) benzo(k)	nene, pyrene, chrysend anthracene, benzo(b) ofluoranthene, benzo( ,2,3-cd)pyrene	fluoranthene,	J	< quantita	tion limit		

Case Number:	31635	Project Number	03-0474	SOW Number		OLC03.2, OLM04.2
Site ID.	Gulf Stat	es Creosoting, Flowd	ood, MS		Date	6/07/03
Affected Samples	Compou	nd or Fraction		Flag <u>Used</u>	Reas	on
6764	chrysene benzo(k)	ene, pyrene, benzo(a, benzo(b)fluoranther fluoranthene, benzo( ,2,3-cd)pyrene	ie,	, <b>J</b>	< quantita	tion limit
6765	phenanth pyrene, b benzo(b)	hyde, acenaphthene, trene, anthracene, flucenzo(a)anthracene, cfluoranthene, benzo(pyrene, indeno(1,2,3	oranthene, hrysene, k)fluoranthene,	J	< quantita	ition limit
6766	benzo(a) indeno(1	hylene, phenanthrene anthracene, chrysene ,2,3-cd)pyrene, a,h)anthracene, benz	, benzo(a)pyrene,	J	< quantita	ation limit
6770, 6787, 6811, 6812, 6813	4,6-dinit	orocyclopentadiene, 2 ro-2-methylphenol, lorobenzidine	,4-dinitrophenol,	J	continuin outlier	g calibration
6770	all comp	ounds		J	extraction outliers	n holding time
6770	4,6-dinit	ro-2-methylphenol		J	surrogate	outlier
6773	carbazol	e		J	< quantita	ation limit
6774	chrysene	nene, pyrene, benzo(a e, benzo(b)fluoranthe ofluoranthene, indene	ne,	J	< quantita	ation limit
6775	benzalde	ehyde, pyrene		J	< quantita	ation limit
6772, 6776, 6777, 6778, 6779, 6780, 6781, 6782	2,4-dini	rophenol		R	poor instr response calibratio	during
6783, 6784, 6785, 6786, 6788, 6789, 6790, 6791, 6792, 6793, 6794, 6795	2,4-dini	trophenol		. J		g calibration RPD outliers

		•	•			•	
Case Number:	31635	Project Number	03-0474	SOW N	Jumber	OLC03.2, OLM04.2	
Site ID.	Gulf States	Creosoting, Flower	ood, MS		Date	6/07/03	
Affected Samples	Compound	or Fraction		Flag <u>Used</u>	Reaso	on	
6772, 6776, 6777, 6778, 6779, 6780, 6781, 6782	4-nitroanili pentachloro	cyclopentadiene, 4 ne, hexachloroben ophenol, indeno(1, a)anthracene	zene,	J	continuing %D outlie	calibration r	
6778	benzo(a)an indeno(1,2,	lene, phenanthrend thracene, benzo(a 3-cd)pyrene, n)anthracene, benz	)pyrene,	J	< quantita	tion limit	
6779		lene, anthracene, o n)anthracene, benz	•	J	< quantita	tion limit	
6783	phenanthre	ne, carbazole		J	< quantita	tion limit	
6784	acenaphthy	lene, carbazole		J	< quantitation limit		
6785, 6786	-	e, 2-methylnaphth ne, dibenzofuran, ophenol	alene,	J	< quantita	tion limit	
6787	all compou	nds		· J	extraction outliers	holding time	
6787	2,4-dimeth dimethylph dinitro-2-n butylbenzy	nenol, 4-methylpho ylphenol, caprolac athalate, diethylpho nethylphenol, di-n- alphthalate, bis(2- alphthalate, di-n-oc	tam, thalate, 4,6-	1	surrogate	outlier	
6788	acenaphthy pentachlor	e, 2-methylnaphthylene, dibenzofura ophenol, carbazole h)anthracene	n,	J ·	< quantita	tion limit	
6789	anthracene	, benzo(k)fluorant	hene	J	< quantita	tion limit	
6791, 6793, 6794, 6795	benzaldeh	yde		<sub>.</sub> J	< quantita	ition limit	

		·	OLC03.2,
Case Number:	31635 Project Number 03-0474	SOW 1	Number OLM04.2
Site ID.	Gulf States Creosoting, Flowood, MS	Date 6/07/03	
Affected Samples	Compound or Fraction	Flag <u>Used</u>	Reason
67.95	hexachlorocyclopentadiene	J	continuing calibration %D outlier
6796, 6797, 6798, 6799, 6800, 6802, 6804, 6805, 6806	2,4-dinitrophenol, 4,6-dinitro-2-methylphenol	J	continuing calibration %D outlier
6801, 6807, 6808, 6809, 6810	butylbenzylphthalate, bis(2-ethylhexyl)phthalate	·· J	continuing calibration %D outlier
6797, 6802, 6804, 6805, 6806, 6809	benzo(b)fluoranthene, benzo(k)fluoranthene	J	isomers could not be resolved by GC
6797	acenaphthylene, anthracene, indeno(1,2,3-cd)pyrene, dibenzo(a,h)anthracene, benzo(g,h,i)perylene	J	< quantitation limit
6802	acenaphthylene, phenanthrene, anthracene, carbazole	J	< quantitation limit
6804	anthracene, fluoranthene, pyrene, chrysene, benzo(a)anthracene, benzo(a)pyrene, dibenzo(a,h)anthracene, benzo(g,h,i)perylene	J	< quantitation limit
6805	naphthalene, 2-methylnaphthalene, 1,1'-biphenyl, acenaphthylene, acenaphthene, dibenzofuran, fluorene	J	< quantitation limit
6806	benzo(b)fluoranthene, benzo(k)fluoranthene	J	< quantitation limit
6809	acenaphthylene, phenanthrene, carbazole, dibenzo(a,h)anthracene, benzo(g,h,i)perylene		< quantitation limit
6811	caprolactam, 1,1-biphenyl, dimethylphthalate, diethylphthalate, 4,6-dinitro-2-methylphenol, di-n-butylphthalate, butylbenzylphthalate, bis(2-ethylhexyl)phthalate, di-n-octylphthalate	J	surrogate outlier
6812	4,6-dinitro-2-methylphenol	J	surrogate outlier

## ORGANIC DATA QUALIFIER REPORT

Case Number:	31635 Project Number 03-0474	SOW N	umber	OLC03.2, OLM04.2
Site ID.	Gulf States Creosoting, Flowood, MS		Date	6/07/03
Affected Samples	Compound or Fraction	Flag <u>Used</u>	Reaso	<u>n</u>
PAHs by GC/MS SIM				
6754	naphthalene, 2-methylnaphthalene, acenaphthene, fluorene, benzo(a)anthracene, chrysene	J	< quantitat	ion limit
6759	naphthalene, 2-methylnaphthalene, acenaphthene, dibenzo(a,h)anthracene	J	< quantitat	ion limit
6775	naphthalene, 2-methylnaphthalene	J	< quantitat	ion limit
6776	naphthalene, 2-methylnaphthalene, acenaphthene, fluorene	<b>J</b>	< quantitat	ion limit
6777	naphthalene, 2-methylnaphthalene	J	< quantitat	ion limit
6792	naphthalene, 2-methylnaphthalene, fluorene	. <b>J</b>	< quantitat	ion limit
6793	naphthalene	J,	< quantitat	ion limit

## ORGANIC DATA QUALIFIER REPORT

Case Number:	31635 Project Number 03-0474	SOW <sup>1</sup>	Number	OLC03.2, OLM04.2
Site ID.	Gulf States Creosoting, Flowood, MS		Date	6/07/03
Affected Samples Pesticides	Compound or Fraction	Flag <u>Used</u>	Reas	on
6755	gamma-BHC (Lindane), Aroclor 1260	J ·	< quantita	ition limit
6757	Aroclor 1260	J	< quantita	ation limit
6759	Endosulfan II	J	< quantita	ation limit
6759	Endosulfan sulfate	N	%D betw	een columns
6760	Endrin, Endrin ketone	J	surrogate	outlier
6762	4,4'-DDE	J	< quantita	ation limit
6762	4,4'-DDE	. <b>N</b>	%D betw	een columns
6764	4,4'-DDE	J	< quantita	ation limit
6764	4,4'-DDE	N	%D betw	een columns
6765	Endosulfan sulfate	J	< quantita	ation limit
6765	Endosulfan sulfate	N	%D betw	een columns
6766	4,4'-DDE	J	< quantita	ation limit
6766	Endrin ketone	J	surrogate	outlier
6767	gamma-BHC (Lindane)	J	< quantit	ation limit
6770, 6787, 6811, 6812, 6813	4,4'-DDD, 4,4'-DDE, 4,4'-DDT, Endrin, Endrin ketone, Endrin aldehyde	J		own (poor erformance)
6773	Endrin, Endrin aldehyde	J.	surrogate	outlier
6778	alpha-BHC	· N	%D betw	een columns
6778	alpha-BHC	J	< quantit	ation limit
6778	alpha-BHC, Endrin ketone	J	surrogate	outlier
6779	4,4'-DDT, Endrin aldehyde	N	%D betw	een columns
6779	4,4'-DDT, Endrin aldehyde	J	surrogate	outlier

## ORGANIC DATA QUALIFIER REPORT

Case Number:	31635	Project Number	03-0474	SOW	Number	OLC03.2, OLM04.2
Site ID.	Gulf State	es Creosoting, Flow	ood, MS		Date	6/07/03
Affected Samples	Compour	nd or Fraction		Flag <u>Used</u>	Reaso	on
6783	4,4'-DDT	, Methoxychlor		N	%D betwe	en columns
6783	4,4'-DDT	, Methoxychlor		J	surrogate o	outlier
6784	4,4'-DDT	, ·		J	surrogate o	outlier
6785	Methoxy	chlor		N	%D betwe	en columns
6785	4,4'-DDT	, Methoxychlor		J	surrogate o	outlier
6786	4,4'-DDT	, endrin ketone		N	%D betwe	en columns
6788	4,4'-DDT			J	surrogate o	outlier
6802	4,4'-DDT	, Methoxychlor, En	drin aldehyde	J	surrogate (	outlier
6802	Methoxy	chlor		N	%D betwee	en columns
6805	4,4'-DD1			J	surrogate	outlier
6809	Endrin al	dehyde		J	surrogate	outlier

6793 FY 2003

Sample

Project: 03-0474

Produced by: Goddard, Denise

Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 15:50 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS08SD / MD No: 1XX3 Org Contractor: LIBRTY D No: 1XX3 Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** RESULTS UNITS **ANALYTE** 14 U UG/KG Dichlorodifluoromethane 14 U UG/KG Dibromochloromethane 14 U UG/KG 1,2-Dibromoethane (EDB) 14 U UG/KG Chloromethane 14 U 14 U UG/KG Vinyl Chloride UG/KG Chlorobenzene 14 U UG/KG Bromomethane 14 U UG/KG Ethyl Benzene 14 U UG/KG Chloroethane 14 U UG/KG **Total Xylenes** Trichlorofluoromethane (Freon 11) 14 U UG/KG 14 U UG/KG Styrene 1,1-Dichloroethene (1,1-Dichloroethylene) 14 U UG/KG 14 U UG/KG Bromoform UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 14 U 14 U UG/KG Isopropylbenzene 29 J UG/KG Acetone - 14 U 1,1,2,2-Tetrachloroethane UG/KG 14 U UG/KG Carbon Disulfide 14 U UG/KG 1,3-Dichlorobenzene 14 U UG/KG Methyl Acetate 14 U UG/KG 1.4-Dichlorobenzene 14 U UG/KG Methylene Chloride 14 U UG/KG 1,2-Dichlorobenzene 14 U UG/KG trans-1,2-Dichloroethene 1,2-Dibromo-3-Chloropropane (DBCP) 14 U UG/KG 14 U UG/KG Methyl T-Butyl Ether (MTBE) 14 U UG/KG 1,2,4-Trichlorobenzene 1.1-Dichloroethane 14 U UG/KG NA UG/KG 1.2.3-Trichlorobenzene 14 U UG/KG cis-1,2-Dichloroethene 19 % Moisture 14 UJ UG/KG Methyl Ethyl Ketone UG/KG Bromochloromethane NA 14 U UG/KG Chloroform 14 U UG/KG 1,1,1-Trichloroethane 14 U · UG/KG Cyclohexane 14 U UG/KG Carbon Tetrachloride 14 U UG/KG Benzene 14 U UG/KG 1,2-Dichloroethane 14 U UG/KG Trichloroethene (Trichloroethylene) 14 U UG/KG Methylcyclohexane 1,2-Dichloropropane 14 U UG/KG Bromodichloromethane UG/KG 14 U 14 U UG/KG cis-1,3-Dichloropropene 14 UJ UG/KG Methyl Isobutyl Ketone 14 U UG/KG Toluene 14 U UG/KG trans-1,3-Dichloropropene UG/KG 14 U 1.1.2-Trichloroethane 14 U UG/KG Tetrachloroethene (Tetrachloroethylene) 14 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Project: 03-0474

Sample 6793 FY 2003

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS08SD / Media: SEDIMENT

MD No: 1XX3

D No: 1XX3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 15:50

Ending:

**RESULTS UNITS** 

ANALYTE

8 J UG/KG UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

6792 FY 2003

Project: 03-0474

Sample

Produced by: Goddard, Denise

Program: SF Id/Station: GS06	ates Creosoting Flo	wood, MS Case No: 31635 MD No: 1XX2 Inorg Contractor: SENTIN			Requestor: Project Leader: BSTRIGGO Beginning: 04/22/2003 15:15 Ending:			
Media: SEDIME	NT	D No: 1XX2	Org Contractor: LI	BRIY'	DATA	REPORTED ON DRY WEIGHT BASIS	·	
RESULTS UNI  13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/	CG Dichlorodifluoromethan CG Chloromethane CG Vinyl Chloride CG Bromomethane CG Chloroethane CG Trichlorofluoromethane CG 1,1-Dichloroethene (1,1 CG 1,2-Trichloro-1,2,2-Tri CG Acetone CG Carbon Disulfide CG Methyl Acetate CG Methyl Acetate CG Methyl Acetate CG Methyl T-Butyl Ether (M CG 1,1-Dichloroethane CG Cis-1,2-Dichloroethene CG Cis-1,2-Dichloroethene	(Freon 11) -Dichloroethylene) fluoroethane (Freon 113) e		RESULTS  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  13 U  NA	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	ANALYTE Dibromochloromethane 1,2-Dibromoethane (EDB) Chlorobenzene Ethyl Benzene Total Xylenes Styrene Bromoform Isopropylbenzene 1,1,2,2-Tetrachloroethane 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dibromo-3-Chloropropane (DBCP) 1,2,4-Trichlorobenzene 1,2,3-Trichlorobenzene % Moisture		
13 UJ UG/ NA UG/ NA UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/ 13 U UG/	GG Bromochloromethane GG Chloroform GG 1,1,1-Trichloroethane GG Cyclohexane GG Carbon Tetrachloride GB Benzene GG 1,2-Dichloroethane GG Trichloroethene (Trichloroethene) GG Methylcyclohexane GG 1,2-Dichloropropane GG Bromodichloromethane GG cis-1,3-Dichloropropene GG Methyl Isobutyl Ketone GG Toluene GG trans-1,3-Dichloroprope 1,1,2-Trichloroethane GG Tetrachloroethene (Tetra	ne						

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

6776

FY 2003

Produced by: Goddard, Denise

Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 12:15 Program: SF Case No: 31635 Endina: Id/Station: GS05SD / MD No: 1XT6 Inorg Contractor: SENTIN Org Contractor: LIBRTY D No: 1XT6 Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS ANALYTE **RESULTS UNITS** ANALYTE **RESULTS UNITS** 29 U UG/KG Dichlorodifluoromethane 29 U UG/KG Dibromochloromethane 29 U 29 U UG/KG Chloromethane UG/KG 1,2-Dibromoethane (EDB) 29 U 29 U UG/KG Vinyl Chloride UG/KG Chlorobenzene 29 UJ UG/KG Bromomethane 29 U UG/KG Ethyl Benzene 29 U UG/KG Chloroethane 29 U UG/KG **Total Xylenes** 29 U UG/KG Trichlorofluoromethane (Freon 11) 29 U UG/KG Styrene 29 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 29 U UG/KG Bromoform 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 29 U UG/KG 29 U UG/KG Isopropylbenzene 51 J UG/KG Acetone 29 U UG/KG 1,1,2,2-Tetrachloroethane 29 U UG/KG Carbon Disulfide 29 U UG/KG 1,3-Dichlorobenzene 29 U UG/KG Methyl Acetate 29 U UG/KG 1,4-Dichlorobenzene 29 U UG/KG Methylene Chloride 29 U UG/KG 1,2-Dichlorobenzene 29 U UG/KG trans-1,2-Dichloroethene 29 U UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 29 U UG/KG Methyl T-Butyl Ether (MTBE) 29 U UG/KG 1,2,4-Trichlorobenzene 29 U UG/KG 1,1-Dichloroethane 1,2,3-Trichlorobenzene NA UG/KG cis-1,2-Dichloroethene 29 U UG/KG 54 % Moisture 29 U UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 29 U UG/KG Chloroform 29 U UG/KG 1.1.1-Trichloroethane 29 U Cyclohexane UG/KG 29 U UG/KG Carbon Tetrachloride 29 U UG/KG Benzene. 29 U UG/KG 1.2-Dichloroethane 29 U UG/KG Trichloroethene (Trichloroethylene) 29 U UG/KG Methylcyclohexane 29 U UG/KG 1,2-Dichloropropane 29 U UG/KG Bromodichloromethane UG/KG 29 U cis-1,3-Dichloropropene 29 U UG/KG Methyl Isobutyl Ketone 29 U UG/KG Toluene 29 U trans-1,3-Dichloropropene UG/KG 29 U UG/KG 1.1.2-Trichloroethane 29 U UG/KG Tetrachloroethene (Tetrachloroethylene) 29 U UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable,

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6776 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

33 J

Id/Station: GS05SD /

MD No: 1XT6

Media: SEDIMENT

Case No: 31635

D No: 1XT6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:15

Ending:

**ANALYTE RESULTS UNITS** 

UG/KG

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Project: 03-0474 Sample 6777 FY 2003 Requestor: Volatiles Scan Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 12:45 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS07SD / MD No: 1XT7 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XT7 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS ANALYTE** 12 U UG/KG Dichlorodifluoromethane 12 U UG/KG Dibromochloromethane 12 U 12 U UG/KG Chloromethane UG/KG 1,2-Dibromoethane (EDB) 12 U Vinyl Chloride 12 U UG/KG UG/KG Chlorobenzene 12 UJ UG/KG Bromomethane 12 U UG/KG Ethyl Benzene 12 U UG/KG Chloroethane 12 U UG/KG **Total Xylenes** 12 U Trichlorofluoromethane (Freon 11) UG/KG 12 U UG/KG Styrene 12 U 1,1-Dichloroethene (1,1-Dichloroethylene) UG/KG 12 U UG/KG **Bromoform** 12 U 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) UG/KG 12 U UG/KG Isopropylbenzene 49 J UG/KG Acetone 12 U UG/KG 1,1,2,2-Tetrachloroethane 12 U UG/KG Carbon Disulfide 12 U. UG/KG 1,3-Dichlorobenzene 12 U UG/KG Methyl Acetate 12 U UG/KG 1,4-Dichlorobenzene 12 U UG/KG Methylene Chloride 12 U UG/KG 1,2-Dichlorobenzene 12 U trans-1,2-Dichloroethene UG/KG 12 U 1,2-Dibromo-3-Chloropropane (DBCP) UG/KG 12 U Methyl T-Butyl Ether (MTBE) UG/KG 12 U UG/KG 1,2,4-Trichlorobenzene 12 U UG/KG 1.1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene cis-1,2-Dichloroethene 12 U UG/KG 18 % Moisture 12 U UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 12 U UG/KG Chloroform 12 U UG/KG 1,1,1-Trichloroethane 12 U UG/KG Cyclohexane 12 U UG/KG Carbon Tetrachloride 12 U UG/KG Benzene 12 U UG/KG 1,2-Dichloroethane Trichloroethene (Trichloroethylene) 12 U UG/KG 12 U Methylcyclohexane UG/KG 1,2-Dichloropropane 12 U UG/KG Bromodichloromethane 12 U UG/KG 12 U UG/KG cis-1,3-Dichloropropene Methyl Isobutyl Ketone 12 U UG/KG 12 U UG/KG Toluene 12 U UG/KG trans-1,3-Dichloropropene 12 U UG/KG 1.1.2-Trichloroethane 12 U UG/KG Tetrachloroethene (Tetrachloroethylene) UG/KG Methyl Butyl Ketone 12 U

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

70 U

70 U

70 U

70 U

70 U

70 U

70 U

70 U

70 U

70 UJ

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

70 UJ UG/KG

Produced by: Goddard, Denise 6775 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 11:35 Program: SF Case No: 31635 Endina: Inorg Contractor: SENTIN Id/Station: GS04SD / MD No: 1XT5 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XT5 DATA REPORTED ON DRY WEIGHT BASIS **ANALYTE** RESULTS UNITS **ANALYTE RESULTS UNITS** 70 U UG/KG 70 U - UG/KG Dichlorodifluoromethane Dibromochloromethane 70 U 70 U UG/KG 1,2-Dibromoethane (EDB) UG/KG Chloromethane 70 U Chlorobenzene 70 U UG/KG Vinyl Chloride UG/KG 70 U 70 UJ UG/KG Bromomethane UG/KG Ethyl Benzene 70 U UG/KG Chloroethane 70 U UG/KG Total Xylenes 70 U UG/KG Trichlorofluoromethane (Freon 11) 70 U UG/KG Styrene 70 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 70 U Bromoform UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 70 U UG/KG 70 U UG/KG Isopropylbenzene 520 J 1.1,2,2-Tetrachloroethane UG/KG Acetone 70 U UG/KG 70 U UG/KG Carbon Disulfide 70 U UG/KG 1,3-Dichlorobenzene 70 U UG/KG Methyl Acetate 70 U UG/KG 1.4-Dichlorobenzene 70 U UG/KG Methylene Chloride 70 U UG/KG 1,2-Dichlorobenzene 70 U UG/KG trans-1.2-Dichloroethene 70 UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 70 U UG/KG Methyl T-Butyl Ether (MTBE) 70 U UG/KG 1,2,4-Trichlorobenzene 70-U UG/KG 1,1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene cis-1,2-Dichloroethene 70 U UG/KG % Moisture 78 88 J UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane UG/KG Chloroform 70 U 70 U UG/KG 1.1.1-Trichloroethane 70 U UG/KG Cyclohexane 70 U UG/KG Carbon Tetrachloride 70 U UG/KG Benzene 70 U UG/KG 1.2-Dichloroethane

Trichloroethene (Trichloroethylene)

Tetrachloroethene (Tetrachloroethylene)

Methylcyclohexane

1,2-Dichloropropane

Bromodichloromethane

cis-1,3-Dichloropropene

trans-1,3-Dichloropropene

Methyl Isobutyl Ketone

1.1.2-Trichloroethane

Methyl Butyl Ketone

Toluene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6775 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS04SD / Media: SEDIMENT

MD No: 1XT5 D No: 1XT5

Case No: 31635

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:35

Endina:

**RESULTS UNITS** 180 J UG/KG

**ANALYTE** 

**UNKNOWN COMPOUND** 

71 NJ UG/KG **ACETALDEHYDE** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6765 FY 2003 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 10:30 Program: SF Case No: 31635 Ending: inorg Contractor: SENTIN Id/Station: GS03SD / MD No: 1XS5 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XS5 DATA REPORTED ON DRY WEIGHT BASIS **ANALYTE** RESULTS UNITS **RESULTS UNITS ANALYTE** 75 U UG/KG Dichlorodifluoromethane 75 U UG/KG Dibromochloromethane Chloromethane 75 U UG/KG 75 U UG/KG 1,2-Dibromoethane (EDB) 75 U UG/KG Vinvl Chloride 75 U UG/KG Chlorobenzene 75 UJ UG/KG Bromomethane 75 U UG/KG Ethyl Benzene 75 U UG/KG Chloroethane 75 U UG/KG **Total Xylenes** 75 U UG/KG Trichlorofluoromethane (Freon 11) 75 U UG/KG Styrene 75 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 75 U Bromoform UG/KG 75 U 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) UG/KG 75 U UG/KG Isopropylbenzene 350 J UG/KG Acetone 75 U UG/KG 1,1,2,2-Tetrachloroethane 51 J UG/KG Carbon Disulfide 75 U UG/KG 1.3-Dichlorobenzene UG/KG Methyl Acetate 75 U 75 U UG/KG 1,4-Dichlorobenzene 75 U UG/KG Methylene Chloride 75 U UG/KG 1.2-Dichlorobenzene 75 U UG/KG trans-1,2-Dichloroethene 75 U UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 75 U UG/KG Methyl T-Butyl Ether (MTBE) 75 U UG/KG 1,2,4-Trichlorobenzene 75 U UG/KG 1.1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene 75 U UG/KG cis-1,2-Dichloroethene 79 % Moisture 75 U UG/KG Methyl Ethyl Ketone Bromochloromethane NA UG/KG 75 U UG/KG Chloroform 75 U UG/KG 1,1,1-Trichloroethane 75 U UG/KG Cyclohexane 75 U UG/KG Carbon Tetrachloride 75 U UG/KG Benzene 75 U UG/KG 1.2-Dichloroethane 75 U UG/KG Trichloroethene (Trichloroethylene) 75 U UG/KG Methylcyclohexane 75 U UG/KG 1,2-Dichloropropane 75 U UG/KG Bromodichloromethane 75 U UG/KG cis-1,3-Dichloropropene Methyl Isobutyl Ketone 75 U UG/KG 75 U UG/KG Toluene 75 U UG/KG trans-1,3-Dichloropropene 75 U UG/KG 1,1,2-Trichloroethane UG/KG Tetrachloroethene (Tetrachloroethylene) 75 U UG/KG 75 U Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6764 FY 2003 **MISCELLANEOUS COMPOUNDS** 

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 10:45

Ending:

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS20SD /

MD No: 1XS4

Inorg Contractor: SENTIN

Media: SEDIMENT

**D No: 1XS4** 

Org Contractor: LIBRTY

**RESULTS UNITS** 42 NJ UG/KG

**ANALYTE** 

**ACETALDEHYDE** 

7 NJ UG/KG 11 NJ UG/KG 9 J UG/KG

PROPANAL, 2-MEHTYL-BUTANAL, 3-METHYL-UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Sample 6764 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 10:45 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS20SD / MD No: 1XS4 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XS4 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS** ANALYTE 11 U UG/KG Dichlorodifluoromethane 11 U UG/KG Dibromochloromethane 11 U UG/KG Chloromethane .11 U UG/KG 1.2-Dibromoethane (EDB) 11 U UG/KG Vinyl Chloride UG/KG 11 U Chlorobenzene 11 U.J UG/KG Bromomethane 11 U UG/KG Ethyl Benzene 11 U UG/KG Chloroethane UG/KG 11 U Total Xylenes 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene 11 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 11 U UG/KG Bromoform 11 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG Isopropylbenzene 230 J UG/KG Acetone 11 U UG/KG 1,1,2,2-Tetrachloroethane 11 U UG/KG Carbon Disulfide UG/KG 11 U 1.3-Dichlorobenzene 11 U UG/KG Methyl Acetate 11 U UG/KG 1,4-Dichlorobenzene 11 U UG/KG Methylene Chloride 11 U UG/KG 1,2-Dichlorobenzene 11 U UG/KG trans-1,2-Dichloroethene 11 UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) UG/KG 11 U Methyl T-Butyl Ether (MTBE) 11 U UG/KG 1.2.4-Trichlorobenzene 11 U UG/KG 1,1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene 11 U UG/KG cis-1,2-Dichloroethene % Moisture 20 23 J UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 11 U UG/KG Chloroform 11 U UG/KG 1,1,1-Trichloroethane UG/KG Cvclohexane 11 U · 11 U UG/KG Carbon Tetrachloride 11 U UG/KG Benzene 11 U UG/KG 1,2-Dichloroethane 11 U UG/KG Trichloroethene (Trichloroethylene) UG/KG 11 U Methylcyclohexane 11 U UG/KG 1,2-Dichloropropane 11 U UG/KG Bromodichloromethane 11 U UG/KG cis-1.3-Dichloropropene 11 UJ UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene 11 U UG/KG trans-1.3-Dichloropropene 11 U UĠ/KG 1.1.2-Trichloroethane 11 U UG/KG Tetrachloroethene (Tetrachloroethylene) 11 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6759 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:50 Program: SF Case No: 31635 Ending: inorg Contractor: SENTIN Id/Station: GS02SD / MD No: 1XR9 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XR9 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS** ANALYTE 34 U UG/KG Dichlorodifluoromethane 34 U UG/KG Dibromochloromethane 34 U 34 U UG/KG 1.2-Dibromoethane (EDB) UG/KG Chloromethane 34 U UG/KG Chlorobenzene 34 U UG/KG Vinyl Chloride 34 U UG/KG Ethyl Benzene 34 UJ UG/KG Bromomethane 34 U UG/KG Total Xylenes 34 U UG/KG Chloroethane 34 U UG/KG Trichlorofluoromethane (Freon 11) 34 U UG/KG Styrene 1.1-Dichloroethene (1.1-Dichloroethylene) 34 U UG/KG Bromoform 34 U UG/KG 34 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 34 U UG/KG Isopropylbenzene 34 U 130 J UG/KG UG/KG 1,1,2,2-Tetrachloroethane 34 U Carbon Disulfide 34 U UG/KG UG/KG 1.3-Dichlorobenzene 34 U UG/KG Methyl Acetate 34 U UG/KG 1.4-Dichlorobenzene 34 U UG/KG Methylene Chloride 34 U UG/KG 1,2-Dichlorobenzene 34 U UG/KG trans-1,2-Dichloroethene 34 UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 34 U UG/KG Methyl T-Butyl Ether (MTBE) 34 U UG/KG 1.2.4-Trichlorobenzene 34 U UG/KG 1,1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene 34 U UG/KG cis-1,2-Dichloroethene % Moisture 52 34 UJ UG/KG Methyl Ethyl Ketone UG/KG Bromochloromethane NA Chloroform 34 U UG/KG 34 U UG/KG 1,1,1-Trichloroethane 34 U UG/KG Cyclohexane UG/KG Carbon Tetrachloride 34 U 34 U UG/KG Benzene UG/KG 1,2-Dichloroethane 34 U 34 U UG/KG Trichloroethene (Trichloroethylene) UG/KG Methylcyclohexane 34 U 1,2-Dichloropropane 34 U UG/KG Bromodichloromethane 34 U UG/KG 34 U UG/KG cis-1,3-Dichloropropene 34 UJ UG/KG Methyl Isobutyl Ketone UG/KG Toluene 34 U

Data Reported as Identified by CLP Lab - IDs Not Verified

34 U

34 U

34 U

34 UJ

UG/KG

UG/KG

UG/KG

UG/KG

trans-1,3-Dichloropropene

Tetrachloroethene (Tetrachloroethylene)

1.1.2-Trichloroethane

Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6754 FY 2003 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 09:00 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS01SD / MD No: 1XR4

Media: SEDIMENT	D No: 1XR4	D No: 1XR4 Org Contractor: LIBRTY			DATA REPORTED ON DRY WEIGHT BASIS			
RESULTS UNITS AN	ALYTE		RESULTS	UNITS	ANALYTE	:		
37 U UG/KG Dic	hlorodifluoromethane		· 37 U	UG/KG	Dibromochloromethane			
37 U UG/KG Chl	oromethane		37 U	UG/KG	1,2-Dibromoethane (EDB)			
37 U UG/KG Vin	yl Chloride	•	37 U	UG/KG	Chlorobenzene			
37 UJ UG/KG Bro	momethane		37 U	UG/KG	Ethyl Benzene	•		
37 U UG/KG ChI	oroethane		37 U	UG/KG	Total Xylenes			
37 U UG/KG Trid	hlorofluoromethane (Freon 11)		37 U	UG/KG	Styrene			
37 U UG/KG 1,1	Dichloroethene (1,1-Dichloroethylene)		37 U	UG/KG	Bromoform			
37 U UG/KG 1,1,	2-Trichloro-1,2,2-Trifluoroethane (Freon 11	3)	37 U	UG/KG .	Isopropylbenzene	•		
94 J UG/KG Ace	tone		37 U	UG/KG	1,1,2,2-Tetrachloroethane			
	bon Disulfide	•	37 U	UG/KG	1,3-Dichlorobenzene			
37 U UG/KG Met	hyl Acetate		37 U	UG/KG	1,4-Dichlorobenzene			
37 U UG/KG Met	hylene Chloride		37 U	UG/KG	1,2-Dichlorobenzene	•		
37 U UG/KG tran	s-1,2-Dichloroethene	•	37 UJ	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)			
	hyl T-Butyl Ether (MTBE)		37 U	UG/KG	1,2,4-Trichlorobenzene	•		
37 U UG/KG 1,1	Dichloroethane		NA	UG/KG	1,2,3-Trichlorobenzene	•		
37 U UG/KG cis-	1,2-Dichloroethene		-55	%	% Moisture	1		
37 UJ: UG/KG Met	hyl Ethyl Ketone				•			
NA UG/KG Bro	mochloromethane							
37 U UG/KG Chl	oroform							
37 U UG/KG 1,1,	1-Trichloroethane				•			
	lohexane	•						
37 U UG/KG Car	bon Tetrachloride							
37 U UG/KG Ben	zene							
	Dichloroethane				·			
37 U UG/KG Tric	hloroethene (Trichloroethylene)				·			
37 U UG/KG Met	hylcyclohexane							
37 U UG/KG 1,2-	Dichloropropane	•				,		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

Bromodichloromethane

cis-1,3-Dichloropropene

trans-1,3-Dichloropropene

Tetrachloroethene (Tetrachloroethylene)

Methyl Isobutyl Ketone

1,1,2-Trichioroethane

Methyl Butyl Ketone

Toluene

37 U

37 U

37 UJ

37 U

37 U

37 U

37 U

37 UJ

UG/KG

UG/KG

UG/KG

UG/KG UG/KG

UG/KG

UG/KG

UG/KG

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample 6770 FY 2003 Project: 03-0474

Chloroform

Cvclohexane

Benzene

Toluene

1.1.1-Trichloroethane

Carbon Tetrachloride

Trichloroethene (Trichloroethylene)

Tetrachloroethene (Tetrachloroethylene)

1.2-Dichloroethane

Methylcyclohexane

1,2-Dichloropropane Bromodichloromethane

cis-1,3-Dichloropropene

trans-1,3-Dichloropropene

Methyl Isobutyl Ketone

1,1,2-Trichloroethane

Methyl Butyl Ketone

Volatiles Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

5:0 U

0.50 U

0.50 U

0.50 U

0.50 U

5.0 U

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

1011000, 1110

Id/Station: GS01GW /

Case No: 31635

D No: 1XT0

MD No: 1X

Media: GROUNDWATER

MD No: 1XTO Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:05

Ending:

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
0.50 U	UG/L	Dichlorodifluoromethane	0.50 U	UG/L	Dibromochloromethane
0.50 U	UG/L	Chloromethane	0.50 U	UG/L	1,2-Dibromoethane (EDB)
.0.50 U	UG/L	Vinyl Chloride	0.50 U	UG/L	Chlorobenzene
0.50 U	UG/L	Bromomethane	0.50 U	UG/L	Ethyl Benzene
0.50 U	UG/L	Chloroethane	0.50 U	UG/L	Total Xylenes
0.50 U	UG/L	Trichlorofluoromethane (Freon 11)	0.50 U	UG/L	Styrene
0.50 U	UG/L	1,1-Dichloroethene (1,1-Dichloroethylene)	0.50 U	UG/L	Bromoform
0.50 U	UG/L	1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	0.50 U	UG/L	isopropylbenzene
5.0 U	UG/L	Acetone	0.50 U	UG/L	1,1,2,2-Tetrachloroethane
0.50 U	UG/L	Carbon Disulfide	0.50 U	UG/L	1,3-Dichlorobenzene
0.50 U	UG/L	Methyl Acetate	0.50 U	UG/L	1,4-Dichlorobenzene
0.50 U	UG/L	Methylene Chloride	0.50 U	UG/L	1,2-Dichlorobenzene
0.50 U	UG/L	trans-1,2-Dichloroethene	0.50 U	UG/L	1,2-Dibromo-3-Chloropropane (DBCP) -
0.50 U	UG/L	Methyl T-Butyl Ether (MTBE)	0.50 U	UG/L	1,2,4-Trichlorobenzene
0.50 U	UG/L	1,1-Dichloroethane	0.50 U	UG/L	1,2,3-Trichlorobenzene
0.50 U	UG/L	cis-1,2-Dichloroethene			
6.1 UJ	UG/L	Methyl Ethyl Ketone			
0.50 U	UG/L	Bromochloromethane			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6787 FY 2003

Project: 03-0474

1.2-Dichloroethane

Methylcyclohexane

1,2-Dichloropropane Bromodichloromethane

cis-1,3-Dichloropropene

trans-1.3-Dichloropropene

Methyl Isobutyl Ketone

1.1.2-Trichloroethane

Methyl Butyl Ketone

Toluene

Trichloroethene (Trichloroethylene)

Tetrachloroethene (Tetrachloroethylene)

**Volatiles Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS02GW /

0.50 U

0.50 U

0.50 U 0.50 U

0.50 U

0.50 U

5.0 U

0.50 U

0.50 U

0.50 U

0.50 U

5.0 U

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

Flowood, MS

MD No: 1XW7

Case No: 31635 Inorg Contractor: SENTIN

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:55

Ending:

	Media: GR	OUNDWA	TER	D No: 1XW7	Org Contractor: A4	•		
	RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE
•	0.50 U	UG/L	Dichlorodifluoromethane			0.50 U	UG/L	Dibromochloromethane
	0.50 U	UG/L	Chloromethane			.0.50 U	UG/L	1,2-Dibromoethane (EDB)
	0.50 U	UG/L	Vinyl Chloride			0.50 U	UG/L	Chlorobenzene
	0.50 U	UG/L	Bromomethane		•	0.50 U	UG/L	Ethyl Benzene
	0.50 U	UG/L	Chloroethane	•		0.50 U	UG/L	Total Xylenes
	0.50 U	UG/L	Trichlorofluoromethane (	(Freon 11)		0.50 U	UG/L	Styrene
	0.50 U	UG/L	1,1-Dichloroethene (1,1-l	Dichloroethylene)		0.50 U	UG/L	Bromoform
	0.50 U	UG/L	1,1,2-Trichloro-1,2,2-Trifl	uoroethane (Freon 113)	•	0.50 U	UG/L	Isopropylbenzene
	5.0 U	UG/L	Acetone		•	0.50 U	UG/Ł	1,1,2,2-Tetrachloroethane
	0.50 U	UG/L	Carbon Disulfide			0.50 Ų	UG/L	1,3-Dichlorobenzene
	0.50 U	UG/L	Methyl Acetate			0.50 U	UG/L	1,4-Dichlorobenzene
	0.50 U	UG/L	· Methylene Chloride			0.50 U	UG/L	1,2-Dichlorobenzene
	0.50 U	UG/L	trans-1,2-Dichloroethene			0.50 U	UG/L	1,2-Dibromo-3-Chloropropane (DBCP)
	0.50 U	UG/L	Methyl T-Butyl Ether (MT	BE)		0.50 U	UG/L	1,2,4-Trichlorobenzene
	0.50 U	UG/L	1,1-Dichloroethane			0.50 U	UG/L	1,2,3-Trichlorobenzene
	0.50 U	UG/L	cis-1,2-Dichloroethene					
	5.0 UJ	UG/L	Methyl Ethyl Ketone					
	0.50 U	UG/L	Bromochloromethane					
	0.50 U	UG/L	Chloroform					
	0.50 U	UG/L	1,1,1-Trichloroethane					
	0.50 U	UG/L	Cyclohexane					
	0.50 U	UG/L	Carbon Tetrachloride	•			•	•
'	0.50 U	UG/L	Benzene		· ·			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6811 FY .2003

Project: 03-0474

Produced by: Goddard, Denise Requestor:

**Volatiles Scan** 

Facility: Gulf States Creosoting

Flowcod, MS

Project Leader: BSTRIGGO

Program: SF

Id/Station: GS03GW /

Case No: 31635

Beginning: 04/23/2003 09:55 Ending:

Media: GROUNDWATER

MD No: 1XZ1 D No: 1XZ1

Inorg Contractor: SENTIN

Org Contractor: A4

RESULTS	UNITS	ANALYTE	RESULTS UNITS	ANALYTE
0.50 U	UG/L	Dichlorodifluoromethane	0.50 U UG/L	Dibromochloromethane
0.50 U	UG/L	Chloromethane	0.50 U UG/L	1,2-Dibromoethane (EDB)
0.50 U	UG/L	Vinyl Chloride	0.50 U UG/L	Chlorobenzene
0.50 U	UG/L	Bromomethane	0.50 U UG/L	Ethyl Benzene
0.50 U	UG/L	Chloroethane	0.50 U UG/L	Total Xylenes
0.50 U	UG/L	Trichlorofluoromethane (Freon 11)	. 0.50 U UG/L	Styrene
0.50 U	UG/L	1,1-Dichloroethene (1,1-Dichloroethylene)	0.50 U UG/L	Bromoform
0.50 U	UG/Ľ	1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	0.50 U UG/L	Isopropylbenzene
5.0 U	UG/L	Acetone	0.50 U UG/L	1,1,2,2-Tetrachloroethane
0.50 U	UG/L	Carbon Disulfide	0.50 U UG/L	1,3-Dichlorobenzene
0.50 U	UG/L	Methyl Acetate	0.50 U UG/L	1,4-Dichlorobenzene
0.50 U	UG/L	Methylene Chloride	0.50 U UG/L	1,2-Dichlorobenzene
0.50 U	UG/L	trans-1,2-Dichloroethene	0.50 U UG/L	1,2-Dibromo-3-Chloropropane (DBCP)
0.50 U	UG/L	Methyl T-Butyl Ether (MTBE)	0.50 U UG/L	1,2,4-Trichlorobenzene
0.50 U	UG/L	1,1-Dichloroethane	0.50 U UG/L	1,2,3-Trichlorobenzene
0.50 U	UG/L	cis-1,2-Dichloroethene		
6.9 UJ	UG/L	Methyl Ethyl Ketone	•	•
0.50 U	UG/L	Bromochloromethane	•	
0.50 U	UG/L	Chloroform		,
0.50 U	UG/L	1,1,1-Trichloroethane		
0.50 U	UG/L	Cyclohexane		
0.50 U	UG/L	Carbon Tetrachloride		
0.50 U	UG/L	Benzene		
0.50 U	UG/L	1,2-Dichloroethane		
0.50 U	UG/L	Trichloroethene (Trichloroethylene)		•
0.50 U	UG/L	Methylcyclohexane		•
0.50 U	UG/L	1,2-Dichloropropane	•	
0.50 U	UG/L	Bromodichloromethane		
0.50 U	UG/L	cis-1,3-Dichloropropene		•
5.0 U	UG/L	Methyl Isobutyl Ketone		
0.50 U	UG/L	Toluene		•
0.50 U	UG/L	trans-1,3-Dichloropropene		
0.50 U	UG/L	1,1,2-Trichloroethane		
0.50 U	UG/L	Tetrachloroethene (Tetrachloroethylene)		
5.0 U	UG/L	Methyl Butyl Ketone	• .	
				·
		1		·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample **6811** 

6811 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

MISCELLANEOUS COMPOUNDS

Requestor:

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO

Program: SF

Case No: 31635

Beginning: 04/23/2003 09:55

Id/Station: GS03GW /

MD No: 1XZ1

Inorg Contractor: SENTIN

Ending:

Media: GROUNDWATER

D No: 1XZ1

Org Contractor: A4

RESULTS UNITS 1.3 NJ UG/L ANALYTE INDANE

Data Reported as Identified by CLP Lab - IDs Not Verified

Project: 03-0474

Volatiles Scan

Facility: Gulf States Creosoting

6812

Flowood, MS

Program: SF

Sample

Case No: 31635

Id/Station: GS04GW /

FY 2003

Media: GROUNDWATER

MD No: 1XZ3 D No: 1XZ3

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 13:20

Ending:

	RESULTS 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U	UNITS UG/L UG/L UG/L UG/L UG/L UG/L UG/L UG/L	ANALYTE Dichlorodifluoromethane Chloromethane Vinyl Chloride Bromomethane Chloroethane Trichlorofluoromethane ( 1,1-Dichloroethene (1,1-L 1,1,2-Trichloro-1,2,2-Trifluoromethane Carbon Disulfide Methyl Acetate	Dichloroethylene)	113)		RESULTS 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U 0.50 U	UNITS UG/L UG/L UG/L UG/L UG/L UG/L UG/L UG/L	ANALYTE Dibromochloromethane 1,2-Dibromoethane (EDB) Chlorobenzene Ethyl Benzene Total Xylenes Styrene Bromoform Isopropylbenzene 1,1,2,2-Tetrachloroethane 1,3-Dichlorobenzene 1,4-Dichlorobenzene	
	0.50 U 0.50 U	UG/L UG/L	Methylene Chloride trans-1,2-Dichloroethene				0.50 U 0.50 U	UG/L UG/L	1,2-Dichlorobenzene 1,2-Dibromo-3-Chloropropane (DBCP)	
	0.50 U	UG/L	Methyl T-Butyl Ether (MTI	3E)			0.50 U	UG/L	1,2,4-Trichlorobenzene	
	0.50 U	UG/L	1,1-Dichloroethane	J <b>L</b> )			0.50 U	UG/L	1,2,3-Trichlorobenzene	·
	0.50 U	UG/L	cis-1,2-Dichloroethene	•			0.00 0	00/2.	1,2,0 111011010001120110	
	7.2 J	UG/L	Methyl Ethyl Ketone							
	0.50 U	UG/L	Bromochloromethane	•						
٠	0.50 U	UG/L	Chloroform							•
	0.50 U	UG/L	1,1,1-Trichloroethane							•
	0.50 U	UG/L	Cyclohexane							
•	0.50 U	UG/L	Carbon Tetrachloride							
	0.50 U	UG/L	Benzene							
	0.50 U	UG/L	1,2-Dichloroethane						•	
	0.50 U	UG/L	Trichloroethene (Trichloro	ethylene)						
	0.50 U 0.50 U	UG/L UG/L	Methylcyclohexane						,	1
	0.50 U	UG/L	1,2-Dichloropropane Bromodichloromethane							
	0.50 U	UG/L	cis-1,3-Dichloropropene			•				
	5.0 U	UG/L	Methyl Isobutyl Ketone						·	
	0.50 U	UG/L	Toluene	•					e.	
	0.50 U	UG/L	trans-1,3-Dichloropropene	·				•		
	0.50 U	UG/L	1,1,2-Trichloroethane					-		
	0.50 U	·UG/L	Tetrachloroethene (Tetrac	hloroethylene)					•	
	5.0 U	UG/L	Methyl Butyl Ketone						·	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6813 FY 2003 Project: 03-0474

**Volatiles Scan** 

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

0.50 U

5.0 U

5.0 U

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L

UG/L UG/L

UG/L

UG/L

UG/L

UG/L

Carbon Tetrachloride

Trichloroethene (Trichloroethylene)

Tetrachloroethene (Tetrachloroethylene)

1.2-Dichloroethane

Methylcyclohexane

1,2-Dichloropropane

Bromodichloromethane

cis-1,3-Dichloropropene Methyl Isobutyl Ketone

trans-1.3-Dichloropropene

1,1,2-Trichloroethane

Methyl Butyl Ketone

Benzene

Toluene

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

riowood, MS

Id/Station: GS05GW /

Case No: 31635

MD No: 1XZ4

D No: 1X74

.

Inorg Contractor: SENTIN

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 11:45

Ending:

Media: GR	OUNDWA	TER :	D No: 1XZ4	Org Contractor: A4		
RESULTS	UNITS	ANALYTE			RESUL'	
0.50 U	UĠ/L	Dichlorodifluoromethane			0.50 เ	
0.50 U	UG/L	Chloromethane		•	0.50 L	
0.50 U	UG/L	Vinyl Chloride			0.50 L	
0.50 U	UG/L	Bromomethane	• •		0.50 L	
0.50 U	UG/L	Chloroethane			0.50 L	
0.50 U	UG/L	Trichlorofluoromethane (F	Freon 11)	•	0.50 L	
0.50 U	UĠ/L	1,1-Dichloroethene (1,1-D	ichloroethylene)		0.50 L	
0.50 U	UG/L	1,1,2-Trichloro-1,2,2-Triflu	oroethane (Freon 113)		0.50 L	
5.0 U	UG/L	Acetone			0.50 L	
0.50 U	UG/L	Carbon Disulfide			0.50 L	
0.50 U	UG/L		•		0.50 L	
	UG/L				0.50 \	
		·	•		· 0.50 L	
		, , ,	BE)		0.50 L	
					0.50 L	
0.50 U	UG/L			٥		
	UG/L	, ,				
	UG/L				•	
0.50 U	UG/L					
0.50 U	UG/L					
0.50 U	UG/L	Cyclohexane				
	RESULTS  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U  0.50 U	RESULTS UNITS  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L	0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U UG/L 0.50 U	RESULTS UNITS  0.50 U UG/L Dichlorodifluoromethane 0.50 U UG/L Chloromethane 0.50 U UG/L Vinyl Chloride 0.50 U UG/L Bromomethane 0.50 U UG/L Chloroethane 0.50 U UG/L Chloroethane 0.50 U UG/L Trichlorofluoromethane (Freon 11) 0.50 U UG/L 1,1-Dichloroethene (1,1-Dichloroethylene) 0.50 U UG/L 1,1-Z-Trichloro-1,2,2-Trifluoroethane (Freon 113) 5.0 U UG/L Acetone 0.50 U UG/L Carbon Disulfide 0.50 U UG/L Methyl Acetate 0.50 U UG/L Methyl Acetate 0.50 U UG/L Methylene Chloride 0.50 U UG/L Methyl T-Butyl Ether (MTBE) 0.50 U UG/L 1,1-Dichloroethane 0.50 U UG/L Cis-1,2-Dichloroethene 0.50 U UG/L Methyl T-Butyl Ether (MTBE) 0.50 U UG/L Cis-1,2-Dichloroethene 0.50 U UG/L Methyl Ethyl Ketone 0.50 U UG/L Bromochloromethane 0.50 U UG/L Chloroform 0.50 U UG/L Chloroform	RESULTS UNITS  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  1,1-Dichloroethane  (Freon 11)  0.50 U UG/L  1,1-Dichloroethene (1,1-Dichloroethylene)  0.50 U UG/L  1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)  5.0 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U UG/L  0.50 U 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	RESULTS	UNITS	ANALYTE
	0.50 U	UG/L	Dibromochloromethane
	0.50 U	UG/L	1,2-Dibromoethane (EDB)
	0.50 U	UG/L	Chlorobenzene
	0.50 U	UG/L	Ethyl Benzene
	0.50 U	UG/L	Total Xylenes
٠	0.50 U	UG/L	Styrene
	0.50 U	UG/L	Bromoform
	0.50 U	UG/L	Isopropylbenzene
	0.50 U	UG/L	1,1,2,2-Tetrachloroethane
	0.50 U	UG/L	1,3-Dichlorobenzene
	0.50 U	UG/L	1,4-Dichlorobenzene
	0.50 U	UG/L	1,2-Dichlorobenzene
	· 0.50 U	UG/L	1,2-Dibromo-3-Chloropropane (DBCP)
	0.50 U	UG/L	1,2,4-Trichlorobenzene
	0.50 U	UG/L	1,2,3-Trichlorobenzene
	•		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

10 U

UG/KG

UG/KG

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Project: 03-0474 Sample 6809 FY 2003 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 09:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS05SS / MD No: 1XY9 Org Contractor: LIBRTY D No: 1XY9 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS R

RESULTS	UNITS	ANALYTE	-		RESULTS	UNITS	ANALYTE	•		
10 U	UG/KG	Dichlorodifluoromethane	<b>.</b>	•,	10 U	UG/KG	Dibromochloromethane			
10 U	UG/KG	Chloromethane	•		10 U	UG/KG	1,2-Dibromoethane (EDB)			
10 Ú	UG/KG	Vinyl Chloride	-		10 U	UG/KG	Chlorobenzene	•		
10 UJ	UG/KG	Bromomethane	<u>.</u>		10 U	UG/KG	Ethyl Benzene			
10 U	UG/KG	Chloroethane		•	10 U	UG/KG	Total Xylenes			
10 U	UG/KG	Trichlorofluoromethane	(Freon 11)		10 U	UG/KG	Styrene			
10 U	UG/KG	1,1-Dichloroethene (1,1			10 U	UG/KG	Bromoform			•
10 U	UG/KG		fluoroethane (Freon 113)		10 U	UG/KG	Isopropylbenzene	•		
91 J	UG/KG	Acetone	· · · · · · · · · · · · · · · · · · ·		10 U	· UG/KG	1,1,2,2-Tetrachloroethane			
10 U	UG/KG	Carbon Disulfide		•	10 U	UG/KG	1,3-Dichlorobenzene			
10 U	UG/KG	Methyl Acetate	{		10 U	UG/KG	1,4-Dichlorobenzene	•		
10 U	UG/KG	Methylene Chloride	J ፟፟፟	•	10 U	UG/KG	1,2-Dichlorobenzene		•	
10 U	UG/KG	trans-1,2-Dichloroethene	<del>9</del>		10 UJ	UG/KG	1,2-Dibromo-3-Chloropropane	(DBCP)		
10 UJ	UG/KG	Methyl T-Butyl Ether (M)	ΓBE)		10 U	UG/KG	1,2,4-Trichlorobenzene			
10 U	UG/KG	1,1-Dichloroethane			NA	UG/KG	1,2,3-Trichlorobenzene			
10 U	UG/KG	cis-1,2-Dichloroethene			15	%	% Moisture		•	
10 UJ	UG/KG	Methyl Ethyl Ketone	,		•					
NA	UG/KG	Bromochloromethane	1							
10 U	UG/KG	Chloroform	į				•			
10 U	UG/KG	1,1,1-Trichloroethane								
10 U	UG/KG	Cyclohexane	!	•						
·10 U	UG/KG	Carbon Tetrachloride								
10 U	UG/KG	Benzene				•				
10 U	UG/KG	1,2-Dichloroethane		•					•	
10 U	UG/KG	Trichloroethene (Trichlor	oethylene)					•		
10 U	UG/KG	Methylcyclohexane				•				
10 U	UG/KG	1,2-Dichloropropane	I							
10 U	UG/KG	Bromodichloromethane								
10 U	UG/KG	cis-1,3-Dichloropropene	1					•		
10 UJ	UG/KG	Methyl Isobutyl Ketone	I	,						
10 U	UG/KG	Toluene	1						•	
10 U	UG/KG	trans-1,3-Dichloropropen	ie .				•			
10 U	UG/KG	1,1,2-Trichloroethane								

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Tetrachloroethene (Tetrachloroethylene)

Methyl Butyl Ketone

UG/KG

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6809 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

7 J

Id/Station: GS05SS /

MD No: 1XY9

Media: SURFACE SOIL (0" - 12")

D No: 1XY9

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:15

Ending:

**RESULTS UNITS** 

**ANALYTE** 

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise

6807 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 08:30 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS04SS / MD No: 1XY7 Org Contractor: LIBRTY D No: 1XY7 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS** ANALYTE **RESULTS UNITS ANALYTE** 12 U UG/KG Dichlorodifluoromethane 12 U UG/KG Dibromochloromethane 12 U UG/KG 1,2-Dibromoethane (EDB) 12 U UG/KG Chloromethane 12 U UG/KG Vinyl Chloride 12 U UG/KG Chlorobenzene Bromomethane 12 U UG/KG Ethyl Benzene 12 UJ UG/KG 12 U UG/KG Chloroethane 12 U UG/KG **Total Xylenes** 12 U UG/KG Trichlorofluoromethane (Freon 11) 12 U UG/KG Styrene 12 U UG/KG 1.1-Dichloroethene (1.1-Dichloroethylene) 12 U UG/KG Bromoform 12 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 12 U · UG/KG Isopropylbenzene 28 J UG/KG Acetone 12 U 1,1,2,2-Tetrachloroethane UG/KG 12 U · UG/KG Carbon Disulfide 12 U 1.3-Dichlorobenzene UG/KG UG/KG Methyl Acetate 12 U 12 U UG/KG 1.4-Dichlorobenzene Methylene Chloride 12 U UG/KG 12 U UG/KG 1.2-Dichlorobenzene 12 U trans-1,2-Dichloroethene UG/KG 12 UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) UG/KG Methyl T-Butyl Ether (MTBE) 12 UJ 12 U UG/KG 1,2,4-Trichlorobenzene 12 U UG/KG 1,1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene 12 U UG/KG cis-1,2-Dichloroethene % Moisture 12 UJ UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 12 U UG/KG Chloroform 1,1,1-Trichloroethane 12 U ·UG/KG 12 U UG/KG Cyclohexane 12 U UG/KG Carbon Tetrachloride Benzene 12 U UG/KG 1.2-Dichloroethane 12 U UG/KG Trichloroethene (Trichloroethylene) 12 U UG/KG Methylcyclohexane 12 U UG/KG 12 U UG/KG 1,2-Dichloropropane 12 U UG/KG Bromodichloromethane UG/KG cis-1.3-Dichloropropene 12 U Methyl Isobutyl Ketone 12 UJ UG/KG UG/KG Toluene 12 U 12 U UG/KG trans-1,3-Dichloropropene 12 U UG/KG 1,1,2-Trichloroethane 12 U UG/KG Tetrachloroethene (Tetrachloroethylene) 12 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6807

FY 2003 Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS04SS /

MD No: 1XY7

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XY7

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:30

Ending:

**RESULTS UNITS** 

**ANALYTE** 

.19 J UG/KG **UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

10 UJ UG/KG

Methyl Butyl Ketone

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Sample 6805 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 08:35 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS07SS / MD No: 1XY5 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XY5 DATA REPORTED ON DRY WEIGHT BASIS

	<del></del>		<del></del>		<del></del>					
RESULTS	UNITS	ANALYTE	:			RE	SULTS	UNITS	ANALYTE	
10 U	UG/KG	Dichlorodifluoromethane	e ·				10 U	UG/KG	Dibromochloromethane	
10 U	UG/KG	Chloromethane	i .				10 U	UG/KG	1,2-Dibromoethane (EDB)	
10 U	UG/KG	Vinyl Chloride					10 U	⊍G/KG	Chlorobenzene	
. 10 UJ	UG/KG	Bromomethane	1		•		10 U	UG/KG	Ethyl Benzene	
10 U	UG/KG	Chloroethane					10 U	UG/KG	Total Xylenes	•
10 U	UG/KG	Trichlorofluoromethane	(Freon 11)				10 U	UG/KG	Styrene	
10 U	UG/KG	1,1-Dichloroethene (1,1					10 U	UG/KG	Bromoform	
10 U	ÚG/KG	1,1,2-Trichloro-1,2,2-Tri		3)			10 U	UG/KG	Isopropylbenzene	
87 J	UG/KG	Acetone					10 U	UG/KG	1,1,2,2-Tetrachloroethane	•
10 U	UG/KG	Carbon Disulfide	1				10 U	UG/KG	1,3-Dichlorobenzene	
10 U	UG/KG	Methyl Acetate					10 U	. UG/KG	1,4-Dichlorobenzene	
10 ⊍	UG/KG	Methylene Chloride	1				10 U	UG/KG	1,2-Dichlorobenzene	
10 U	UG/KG	trans-1,2-Dichloroethen					10 UJ	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)	•
10 UJ	UG/KG	Methyl T-Butyl Ether (M	TBE)				10 U	UG/KG	1,2,4-Trichlorobenzene	
10 U	UG/KG	1,1-Dichloroethane	İ.				NA	UG/KG	1,2,3-Trichlorobenzene	
10 U	UG/KG	cis-1,2-Dichloroethene					15	%	% Moisture	
10 UJ	UG/KG	Methyl Ethyl Ketone	Ì							
NA	UG/KĢ	Bromochloromethane	1							
10 Ų	UG/KG	Chloroform	! !							
10 Ü	UG/KG	1,1,1-Trichloroethane								•
10 U	UG/KG	Cyclohexane	i !			•			·	
10 U	UG/KG	Carbon Tetrachloride	į.	•						
10 U	UG/KG	Benzene	•						•	•
10 U	UG/KG	1,2-Dichloroethane				•				
10 U	UG/KG	Trichloroethene (Trichlor	pethylene)			•				
10 U	UG/KG	Methylcyclohexane			•					
10 U	UG/KG	1,2-Dichloropropane							•	
10 U	UG/KG	Bromodichloromethane	i .		•	•				
10 U	UG/KG	cis-1,3-Dichloropropene	; :							
10 UJ	UG/KG	Methyl Isobutyl Ketone	, , , , , , , , , , , , , , , , , , ,		•					
10 U	UG/KG	Toluene	į ·		-					
10 U	UG/KG	trans-1,3-Dichloroproper	ie						•	•
10 U	UG/KG	1,1,2-Trichloroethane					*		,	
10 U	UG/KG	Tetrachloroethene (Tetra	ichloroethylene)							

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Poported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6805 FY 2003 Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS07SS /

Media: SURFACE SOIL (0" - 12")

MD No: 1XY5

D No: 1XY5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:35

Ending:

**ANALYTE RESULTS UNITS** 

10 J UG/KG

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyze analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise 6802 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 08:10 Program: SF Case No: 31635 Endina: Inorg Contractor: SENTIN Id/Station: GS08SS / MD No: 1XY3 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XY3 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS** ANALYTE 11 U UG/KG Dichlorodifluoromethane 11 U UG/KG Dibromochloromethane 11 U UG/KG Chloromethane 11 U UG/KG 1,2-Dibromoethane (EDB) 11 U UG/KG 11 U UG/KG Vinyl Chloride Chlorobenzene 11 UJ UG/KG 11 U UG/KG Ethyl Benzene Bromomethane UG/KG Chloroethane 11 U UG/KG **Total Xylenes** 11 U 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene 1,1-Dichloroethene (1,1-Dichloroethylene) Bromoform 11 U UG/KG 11 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG 11 U UG/KG Isopropylbenzene 74 J UG/KG Acetone 11 U UG/KG 1,1,2,2-Tetrachloroethane 11 U Carbon Disulfide UG/KG 11 U UG/KG 1,3-Dichlorobenzene 11 U UG/KG Methyl Acetate 11 U UG/KG 1.4-Dichlorobenzene Methylene Chloride 11 U UG/KG 11 U UG/KG 1,2-Dichlorobenzene 11 U UG/KG trans-1,2-Dichloroethene 11 UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 11 UJ UG/KG Methyl T-Butyl Ether (MTBE) UG/KG 1,2,4-Trichlorobenzene 11 U 11 U UG/KG 1.1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene 11 U UG/KG cis-1,2-Dichloroethene % Moisture 11 UJ UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane Chloroform 11 U UG/KG 11 U UG/KG 1,1,1-Trichloroethane 11 U UG/KG Cyclohexane 11 U UG/KG Carbon Tetrachloride 11 U UG/KG Benzene 1,2-Dichloroethane UG/KG 11 U 11 U UG/KG Trichloroethene (Trichloroethylene) 11 U UG/KG Methylcyclohexane 11 U UG/KG 1,2-Dichloropropane Bromodichloromethane 11 U UG/KG 11 U UG/KG cis-1.3-Dichloropropene 11 UJ UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene 11 U UG/KG trans-1,3-Dichloropropene UG/KG 1,1,2-Trichloroethane 11 U 11 U. UG/KG Tetrachloroethene (Tetrachloroethylene) 11 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be less than the reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Project: 03-0474

Produced by: Goddard, Denise

6802 FY 2003

Requestor:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:10

Program: SF

18 J

Case No: 31635

Endina:

Id/Station: GS08SS /

MD No: 1XY3 D No: 1XY3

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

UG/KG

Org Contractor: LIBRTY

**RESULTS UNITS** 

**ANALYTE** 

2 UNKNOWN COMPOUNDS

Cyanide Analysis Not Requested

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6800 FY **2003** Project: 03-0474

**Volatiles Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS03SS /

Case No: 31635

MD No: 1XY1 D No: 1XY1

Media: SURFACE SOIL (0" - 12")

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 07:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
13 U	UG/KG	Dichlorodifluoromethane	13 U	UG/KG	Dibromochloromethane
13 U	UG/KG	Chloromethane	13 U	UG/KG	1,2-Dibromoethane (EDB)
13 U	UG/KG	Vinyl Chloride	13 U	UG/KG	Chlorobenzene
13 ÜJ	UG/KG	Bromomethane	13 U	UG/KG	Ethyl Benzene
13 U	UG/KG	Chloroethane	13 U	UG/KG	Total Xylenes
13 U	UG/KG	Trichlorofluoromethane (Freon 11)	13 U	UG/KG	Styrene
13 U	UG/KG	1,1-Dichloroethene (1,1-Dichloroethylene)	13 U	UG/KG	Bromoform
13 U	UG/KG	1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	13 U	UG/KG	Isopropylbenzene
25 J	UG/KG	Acetone	13 U	UG/KG	1,1,2,2-Tetrachloroethane
13 U	UG/KG	Carbon Disulfide	13 U	UG/KG	1,3-Dichlorobenzene
13 U	UG/KG	Methyl Acetate	13 U	UG/KG	1,4-Dichlorobenzene
13 UJ	UG/KG	Methylene Chloride	13 U	UG/KG	1,2-Dichlorobenzene
13 U	UG/KG	trans-1,2-Dichloroethene	13 UJ	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)
13 U	UG/KG	Methyl T-Butyl Ether (MTBE)	13 U	UG/KG	1,2,4-Trichlorobenzene
13 U	UG/KG	1,1-Dichloroethane	· NA	UG/KG	1,2,3-Trichlorobenzene
13 U	UG/KG	cis-1,2-Dichloroethene	24	%	% Moisture
13 UJ	UG/KG	Methyl Ethyl Ketone			
NA	UG/KG	Bromochloromethane			,
13 U	UG/KG	Chloroform			
. 13 U	UG/KG	1,1,1-Trichloroethane			
13 U	UG/KG	Cyclohexane			•
13 U	UG/KG	Carbon Tetrachloride			•
13 U	UG/KG	Benzene	٠		
13 U	UG/KG	1,2-Dichloroethane			•
13 U	UG/KG	Trichloroethene (Trichloroethylene)			
13 U	UG/KG	Methylcyclohexane			
13 U	UG/KG	1,2-Dichloropropane	•	•	•
13 U	UG/KG	Bromodichloromethane			
13 U	UG/KG	cis-1,3-Dichloropropene		•	
13 UJ	UG/KG	Methyl Isobutyl Ketone			
13 U	UG/KG	Toluene	-		
13 U	UG/KG	trans-1,3-Dichloropropene			
	UG/KG	1,1,2-Trichloroethane			
13 U	UG/KG	Tetrachloroethene (Tetrachloroethylene)			
13 UJ	UG/KG	Methyl Butyl Ketone			•
					•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6800

FY **2003** 

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

6 J

Id/Station: GS03SS /

Media: SURFACE SOIL (0" - 12")

MD No: 1XY1

D No: 1XY1

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 07:40

Ending:

**RESULTS UNITS ANALYTE** 

UG/KG

UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Sample FY 2003 Project: 03-0474 6798 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/23/2003 07:40 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS12SS / MD No: 1XX9 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX9 DATA REPORTED ON DRY WEIGHT BASIS

					TIEL CITTED CITTETT INCIGNIT BACK
RESULTS	UNITS	ANALYTE	RESULTS	UNITS.	ANALYTE
10 U	UG/KG	Dichlorodifluoromethane	10 U	UG/KG	Dibromochloromethane
10 U	UG/KG	Chloromethane	10 U	UG/KG	1,2-Dibromoethane (EDB)
10 Ü	UG/KG	Vinyl Chloride	10 U	UG/KG	Chlorobenzene
10 U	UG/KG	Bromomethane	10 U	UG/KG	Ethyl Benzene
10 U	UG/KG	Chloroethane	- 10 U	UG/KG	Total Xylenes
10 U	UG/KG	Trichlorofluoromethane (Freon 11)	10 U	UG/KG	Styrene
10 U	UG/KG	1,1-Dichloroethene (1,1-Dichloroethylene)	· 10 U	UG/KG	Bromoform
10 U	UG/KG	1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	10 U	UG/KG	isopropylbenzene
210 UJ	UG/KG	Acetone	10 U	UG/KG	1,1,2,2-Tetrachloroethane
10 U	UG/KG	Carbon Disulfide		UG/KG	1,3-Dichlorobenzene
10 U	UG/KG	Methyl Acetate		UG/KG	1,4-Dichlorobenzene
10 U	UG/KG	Methylene Chloride	10 U	UG/KG	1,2-Dichlorobenzene
10 U	UG/KG	trans-1,2-Dichloroethene		UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)
10 U	UG/KG	Methyl T-Butyl Ether (MTBE)		UG/KG	1,2,4-Trichlorobenzene
10 U	UG/KG	1,1-Dichloroethane		UG/KG	1,2,3-Trichlorobenzene
10 U	UG/KG	cis-1,2-Dichloroethene	17	%	% Moisture
14 J	UG/KG UG/KG	Methyl Ethyl Ketone			
NA 10 II	UG/KG	Bromochloromethane			
10 U 10 U	UG/KG	Chloroform			•
10 U	UG/KG	1,1,1-Trichloroethane Cyclohexane			
10 U	UG/KG	Carbon Tetrachloride			
10 U	UG/KG	Benzene			
10 U	UG/KG	1,2-Dichloroethane			
10 U	UG/KG	Trichloroethene (Trichloroethylene)			•
10 U	UG/KG	Methylcyclohexane			
10 U	UG/KG	1,2-Dichloropropane			•
10 U	UG/KG	Bromodichloromethane	·		
10 U	UG/KG	cis-1,3-Dichloropropene			
10 UJ	UG/KG	Methyl Isobutyl Ketone			
10 U	UG/KG	Toluene			
	UG/KG	trans-1,3-Dichloropropene			
	UG/KG	`1,1,2-Trichloroethane	•		
	UG/KG	Tetrachloroethene (Tetrachloroethylene)			
10 UJ	UG/KG	Methyl Butyl Ketone			
					•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6798 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS12SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

MD No: 1XX9

D No: 1XX9

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 07:40

Ending:

**RESULTS UNITS ANALYTE** 5 NJ UG/KG BUTANAL

18 NJ UG/KG HEXANAL 30 J UG/KG

3 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise 6797 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:35 Program: SF Case No: 31635 Endina: Inorg Contractor: SENTIN ld/Station: GS24SS / MD No: 1XX7 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX7 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS RESULTS UNITS ANALYTE** ANALYTE 25 U UG/KG Dichlorodifluoromethane 25 U UG/KG Dibromochloromethane 25 U 25 U UG/KG Chloromethane UG/KG 1.2-Dibromoethane (EDB) 25 U UG/KG Vinvl Chloride 25 U UG/KG Chlorobenzene 25 U 25 U UG/KG Bromomethane UG/KG Ethyl Benzene 25 U UG/KG Chloroethane 25 U UG/KG Total Xylenes 25 U UG/KG Trichlorofluoromethane (Freon 11) 25 U UG/KG Styrene 25 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 25 U UG/KG Bromoform 1.1.2-Trichloro-1.2,2-Trifluoroethane (Freon 113) 25 U UG/KG 25 U UG/KG Isopropylbenzene 220 J UG/KG Acetone 25 U UG/KG 1,1,2,2-Tetrachloroethane

25 U UG/KG Carbon Disulfide 25 U UG/KG Methyl Acetate 25 U UG/KG Methylene Chloride trans-1,2-Dichloroethene 25 U UG/KG 25 U UG/KG Methyl T-Butyl Ether (MTBE) 25 U UG/KG 1,1-Dichloroethane 25 U cis-1,2-Dichloroethene UG/KG 25 UJ UG/KG Methyl Ethyl Ketone Bromochloromethane NA UG/KG 25 U UG/KG Chloroform 25 U UG/KG 1,1,1-Trichloroethane 25 U Cyclohexane UG/KG 25 U UG/KG Carbon Tetrachloride 25 U UG/KG Benzene 25 U UG/KG 1,2-Dichloroethane 25 U UG/KG Trichloroethene (Trichloroethylene) 25 U Methylcyclohexane UG/KG 1.2-Dichloropropane 25 U UG/KG 25 U UG/KG Bromodichloromethane 25 U UG/KG cis-1.3-Dichloropropene 25 UJ UG/KG · Methyl Isobutyl Ketone 25 U UG/KG Toluene 25 U UG/KG trans-1,3-Dichloropropene 25 U 'UG/KG 1.1.2-Trichloroethane 25 U UG/KG Tetrachloroethene (Tetrachloroethylene) 25 UJ UG/KG Methyl Butyl Ketone

25 U UG/KG 1,3-Dichlorobenzene 25 U UG/KG 1.4-Dichlorobenzene 25 U UG/KG 1.2-Dichlorobenzene 25 U . · · UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 25 U UG/KG 1,2,4-Trichlorobenzene NA UG/KG 1,2,3-Trichlorobenzene % Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

MISCELLANEOUS COMPOUNDS

Project: 03-0474

6797 FY 2003

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS24SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

D No: 1XX7

MD No: 1XX7

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:35

Ending:

**RESULTS UNITS ANALYTE** 

150 J UG/KG 2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Sample 6795 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:40 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN ld/Station: GS06SS / MD No: 1XX5 Org Contractor: LIBRTY D No: 1XX5 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS** ANALYTE **RESULTS UNITS ANALYTE** 13 U UG/KG Dichlorodifluoromethane Dibromochloromethane 13 U UG/KG 13 U UG/KG 13 U 1.2-Dibromoethane (EDB) **UG/KG** Chloromethane 13 U UG/KG Vinyl Chloride 13 U UG/KG Chlorobenzene 13 U 13 U UG/KG Bromomethane UG/KG Ethyl Benzene Chloroethane 13 U UG/KG Total Xylenes 13 U UG/KG Trichlorofluoromethane (Freon 11) 13 U Styrene 13 U UG/KG UG/KG 13 U 1,1-Dichloroethene (1,1-Dichloroethylene) 13 U UG/KG Bromoform UG/KG 13 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 13 U UG/KG Isopropylbenzene 82 J UG/KG Acetone 13 U UG/KG 1,1,2,2-Tetrachloroethane 13 U UG/KG Carbon Disulfide 13 U UG/KG 1.3-Dichlorobenzene 13 U UG/KG Methyl Acetate 13 U UG/KG 1.4-Dichlorobenzene 13 U UG/KG Methylene Chloride 13 U UG/KG 1,2-Dichlorobenzene 13 U UG/KG trans-1.2-Dichloroethene 13 U UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 13 U UG/KG Methyl T-Butyl Ether (MTBE) 13 U 1,2,4-Trichlorobenzene UG/KG 13 U UG/KG 1,1-Dichloroethane 1.2.3-Trichlorobenzene NA UG/KG 13 U UG/KG cis-1,2-Dichloroethene 26 % Moisture 13 UJ UG/KG Methyl Ethyl Ketone Bromochloromethane NA UG/KG UG/KG Chloroform 13 U 1,1,1-Trichloroethane 13 U UG/KG 13 U UG/KG Cyclohexane 13 U UG/KG Carbon Tetrachloride UG/KG Benzene 13 U 1.2-Dichloroethane 13 U UG/KG Trichloroethene (Trichloroethylene) 13 U UG/KG Methylcyclohexane 13 U ·UG/KG 13 U UG/KG 1,2-Dichloropropane Bromodichloromethane 13 U UG/KG 13 U UG/KG cis-1,3-Dichloropropene 13 UJ UG/KG Methyl Isobutyl Ketone 13 U UG/KG Toluene 13 U UG/KG trans-1,3-Dichloropropene 13 U UG/KG 1.1.2-Trichloroethane 13 U UG/KG Tetrachloroethene (Tetrachloroethylene) 13 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. | L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6795 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 16:40

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS06SS /

Case No: 31635

MD No: 1XX5

Media: SURFACE SOIL (0" - 12")

D No: 1XX5

Inorg Contractor: SENTIN Org Contractor: LIBRTY

**RESULTS UNITS** 20 J

UG/KG

**ANALYTE** 

2 UNKNOWN COMPOUNDS

64 NJ UG/KG **HEXANAL** 

Data Reported as Identified by CLP Lab - IDs Not Verified

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise

Sample 6794 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:42 Program: SF Case No: 31635 Ending: ld/Station: GS22SS / Inorg Contractor: SENTIN MD No: 1XX4 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX4. DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS** ANALYTE 11 U UG/KG Dichlorodifluoromethane 11 U UG/KG Dibromochloromethane 11 U UG/KG 1,2-Dibromoethane (EDB) 11 U UG/KG Chloromethane 11 U UG/KG Vinyl Chloride 11 U UG/KG Chlorobenzene UG/KG Bromomethane UG/KG Ethyl Benzene 11 U 11 U 11 U UG/KG Chloroethane 11 U UG/KG **Total Xylenes** 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene 11 U UG/KG 1.1-Dichloroethene (1.1-Dichloroethylene) 11 U UG/KG Bromoform 11 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) . 11 U UG/KG Isopropylbenzene 110 J UG/KG Acetone UG/KG 1,1,2,2-Tetrachloroethane 11 U 11 U UG/KG Carbon Disulfide 11 U UG/KG 1,3-Dichlorobenzene 11 U UG/KG Methyl Acetate 1.4-Dichlorobenzene 11 U UG/KG 11 U UG/KG Methylene Chloride 1,2-Dichlorobenzene 11 U UG/KG trans-1,2-Dichloroethene 11 U UG/KG - 11 U 1,2-Dibromo-3-Chloropropane (DBCP) UG/KG 11 U Methyl T-Butyl Ether (MTBE) UG/KG 11 U 1.2.4-Trichlorobenzene UG/KG 11 U UG/KG 1.1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene 11 U cis-1,2-Dichloroethene UG/KG 13 % Moisture Methyl Ethyl Ketone 11 UJ UG/KG Bromochloromethane NA UG/KG 11 U UG/KG Chloroform 11 U UG/KG 1,1,1-Trichloroethane 11 U UG/KG Cyclohexane 11 U UG/KG Carbon Tetrachloride 11 U UG/KG Benzene. 11 U UG/KG 1.2-Dichloroethane Trichloroethene (Trichloroethylene) 11 U UG/KG 11 U UG/KG Methylcyclohexane 11 U UG/KG 1,2-Dichloropropane 11 U UG/KG Bromodichloromethane 11 U UG/KG cis-1,3-Dichloropropene 11 UJ UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene 11 U UG/KG trans-1,3-Dichloropropene 11 U 1.1.2-Trichloroethane UG/KG 11 U UG/KG Tetrachloroethene (Tetrachloroethylene) 11 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6794 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

6 J

Id/Station: GS22SS /

Media: SURFACE SOIL (0" - 12")

Case No: 31635 MD No: 1XX4 D No: 1XX4

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:42

Ending:

**RESULTS UNITS** ANALYTE

UG/KG

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

13 UJ

UG/KG

UG/KG

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Project: 03-0474 Sample 6790 FY 2003 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:15 Program: SF Case No: 31635 Ending: Id/Station: GS09SS / MD No: 1XX0 Inorg Contractor: SENTIN Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX0 DATA REPORTED ON DRY WEIGHT BASIS

			1			
RI	ESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
	13 U	UG/KG	Dichlorodifluoromethane	13 U	UG/KG	Dibromochloromethane
	13 U ʻ	UG/KG	Chloromethane	.13 U	UG/KG	1,2-Dibromoethane (EDB)
	13 U	UG/KG -	Vinyl Chloride	13 U	UG/KG	Chlorobenzene
	13 UJ	UG/KG	Bromomethane	13 U	UG/KG	Ethyl Benzene
	13 U	UG/KG	Chloroethane	13 U	UG/KG	Total Xylenes
	13 U	UG/KG	Trichlorofluoromethane (Freon 11)	13 U	UG/KG	Styrene
	13·U	UG/KG	1,1-Dichloroethene (1,1-Dichloroethylene)	13 U	UG/KG	Bromoform
	13 U	UG/KG	1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	13 U	UG/KG	Isopropylbenzene
	13 UJ	UG/KG	Acetone	13 U	UG/KG	1,1,2,2-Tetrachloroethane
	13 U	UG/KG	Carbon Disulfide	13 U	UG/KG	1,3-Dichlorobenzene
	13 U	UG/KG	Methyl Acetate	13 U	UG/KG	1,4-Dichlorobenzene
	13 U	UG/KG	Methylene Chloride	13 U	UG/KG.	1,2-Dichlorobenzene
	13 U	UG/KG	trans-1,2-Dichloroethene	13 U	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)
	13 UJ	UG/KG	Methyl T-Butyl Ether (MTBE)	13 U	UG/KG	1,2,4-Trichlorobenzene
	13 U	UG/KG	1,1-Dichloroethane	NA	UG/KG	1,2,3-Trichlorobenzene
	13 U	UG/KG	cis-1,2-Dichloroethene	26	%	% Moisture
	13 UJ	UG/KG	Methyl Ethyl Ketone	~		
	NA	UG/KG	Bromochloromethane			
	13 U	UG/KG	Chloroform			
	13 U	UG/KG	1,1,1-Trichloroethane	•		•
	13 U	UG/KG	Cyclohexane			•
	13 U	UG/KG	Carbon Tetrachloride			
	13 U	UG/KG	Benzene			
	13 U	UG/KG	1,2-Dichloroethane			
	13 U	UG/KG	Trichloroethene (Trichloroethylene)			
	13 U	UG/KG	Methylcyclohexane			
	13 U	UG/KG	1,2-Dichloropropane			
	13 U	UG/KG	Bromodichloromethane			•
	13 U	UG/KG	cis-1,3-Dichloropropene			
	13 UJ	UG/KG	Methyl Isobutyl Ketone			•
	13 U	UG/KG	Toluene			
	13 U	UG/KG	trans-1,3-Dichloropropene			
	13 U	UG/KG	1,1,2-Trichloroethane			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Tetrachloroethene (Tetrachloroethylene)

Methyl Butyl Ketone

UG/KG

11 UJ UG/KG

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Sample 6788 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:00 Program: SF Case No: 31635 Ending: -Id/Station: GS11SS / Inorg Contractor: SENTIN MD No: 1XW8 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XW8 DATA REPORTED ON DRY WEIGHT BASIS

		- (- ,	F 7			UNIN	THE OTTED ON BITT WEIGHT BROKE	
RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE	
11 U	UG/KG	Dichlorodifluoromethane			11 U	UG/KG	Dibromochloromethane	•
11 U	UG/KG	Chloromethane	·		11 U .	UG/KG	1,2-Dibromoethane (EDB)	
. 11 U	UG/KG	Vinyl Chloride			11 U	UG/KG	Chlorobenzene	
· 11 U	UG/KG	Bromomethane			11 U	UG/KG	Ethyl Benzene	•
11 U	UG/KG	Chloroethane			11 Ù	UG/KG	Total Xylenes	
11 U	UG/KG	Trichlorofluoromethane	(Freon 11)	,	11 U	UG/KG	Styrene	
11 U	UG/KG	1,1-Dichloroethene (1,1-	Dichloroethylene)		11 U	UG/KG	Bromoform	
11 U	UG/KG	1,1,2-Trichloro-1,2,2-Trif	luoroethane (Freon 113)		11 U	UG/KG	Isopropylbenzene	
220 J	UG/KG	Acetone			11 U	UG/KG	1,1,2,2-Tetrachloroethane	
· 11 U	UG/KG	Carbon Disulfide			11 U	UG/KG	1,3-Dichlorobenzene	
11 U	UG/KG	Methyl Acetate			11 U	UG/KG	1,4-Dichlorobenzene	
11 U	UG/KG	Methylene Chloride			11 U	UG/KG	1,2-Dichlorobenzene	
11 U	UG/KG	trans-1,2-Dichloroethene			11 U	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)	
11 U	UG/KG	Methyl T-Butyl Ether (MT	BE)		11 U	UG/KG	1,2,4-Trichlorobenzene	
11 U	UG/KG	1,1-Dichloroethane			NA	UG/KG	1,2,3-Trichlorobenzene	
11 U	UG/KG	cis-1,2-Dichloroethene			· 7	%	% Moisture	
21 J	UG/KG	Methyl Ethyl Ketone						
NA	UG/KG	Bromochloromethane						•
11 U	. UG/KG	Chloroform ·					·	
11 U	UG/KG	1,1,1-Trichloroethane						
11 U	UG/KG	Cyclohexane						
11 U	UG/KG	Carbon Tetrachloride			•			
11 U	UG/KG	Benzene						
. 11 U	UG/KG	1,2-Dichloroethane				-		
11 U	UG/KG	Trichloroethene (Trichlor	oethylene)					
11 U	UG/KG	Methylcyclohexane	•				•	
11 U	UG/KG	1,2-Dichloropropane						
11 U	UG/KG	Bromodichloromethane	-		•			
11 U	UG/KG	cis-1,3-Dichloropropene	•					
11 ÚJ	UG/KG	Methyl Isobutyl Ketone					·	
11 U	UG/KG	Toluene		•			•	
11 U	UG/KG	trans-1,3-Dichloroproper	ıe				•	
11 U	UG/KG	1,1,2-Trichloroethane						

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Tetrachloroethene (Tetrachloroethylene)

Methyl Butyl Ketone

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Project: 03-0474

6788 FY 2003

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS11SS /

MD No: 1XW8

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XW8

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:00

Ending:

RESULTS UNITS **ANALYTE** 35 J UG/KG

3 UNKNOWN COMPOUNDS

37 J UG/KG **HEXANAL** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

UG/KG

Methyl Butyl Ketone

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample **6785** Requestor: Volatiles Scan Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 15:05 Program: SF Case No: 31635 Endina: Inorg Contractor: SENTIN Id/Station: GS10SS / MD No: 1XW5 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XW5 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** RESULTS UNITS ANALYTE 11 U UG/KG Dichlorodifluoromethane 11 U UG/KG Dibromochloromethane 11 U UG/KG 11 U UG/KG Chloromethane 1,2-Dibromoethane (EDB) 11 U UG/KG Vinyl Chloride 11 U UG/KG Chlorobenzene 11 UJ Bromomethane 11 U UG/KG Ethyl Benzene UG/KG 11 U UG/KG Chloroethane 11 U UG/KG Total Xylenes 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene 11 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 11 U UG/KG Bromoform 11 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG Isopropylbenzene 84 J UG/KG Acetone 11 U UG/KG 1,1,2,2-Tetrachloroethane 11 U Carbon Disulfide UG/KG 11 U UG/KG 1,3-Dichlorobenzene 11 U Methyl Acetate UG/KG `11 U UG/KG 1,4-Dichlorobenzene Methylene Chloride 11 U UG/KG 11 U UG/KG 1,2-Dichlorobenzene trans-1,2-Dichloroethene 11 U 11 U UG/KG UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 11 U UG/KG Methyl T-Butyl Ether (MTBE) 11 U UG/KG 1,2,4-Trichlorobenzene 11 U UG/KG 1,1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene 11 U UG/KG cis-1,2-Dichloroethene % Moisture 11 U UG/KG Methyl Ethyl Ketone Bromochloromethane NA UG/KG 11 U UG/KG Chloroform 11 U UG/KG 1,1,1-Trichloroethane 11 U UG/KG. Cyclohexane 11 U UG/KG Carbon Tetrachloride 11 U UG/KG Benzene 11 U UG/KG 1,2-Dichloroethane 11 U UG/KG Trichloroethene (Trichloroethylene) 11 U Methylcyclohexane UG/KG 11 U UG/KG 1,2-Dichloropropane UG/KG Bromodichloromethane 11 U 11 U UG/KG cis-1,3-Dichloropropene 11 U UG/KG Methyl Isobutyl Ketone 11 U Toluene UG/KG 11 U UG/KG trans-1.3-Dichloropropene 11 U UG/KG 1.1.2-Trichloroethane 11 U UG/KG Tetrachloroethene (Tetrachloroethylene)

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Project: 03-0474

6785 FY 2003

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

29 J

Sample

Case No: 31635

Id/Station: GS10SS / Media: SURFACE SOIL (0" - 12")

UG/KG

MD No: 1XW5

D No: 1XW5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 15:05

Ending:

**RESULTS UNITS** 

ANALYTE

3 UNKNOWN COMPOUNDS

45 NJ UG/KG

**HEXANAL** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

11 U 11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U 11 U UG/KG

UG/KG

UG/KG

UG/KG

.UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG UG/KG

UG/KG

Cyclohexane

Benzene

Toluene

Carbon Tetrachloride

Trichloroethene (Trichloroethylene)

Tetrachioroethene (Tetrachioroethylene)

1.2-Dichloroethane

Methylcyclohexane

1.2-Dichloropropane

Bromodichloromethane

cis-1,3-Dichloropropene

trans-1,3-Dichloropropene 1,1,2-Trichloroethane

Methyl Isobutyl Ketone

Methyl Butyl Ketone

**EPA - REGION IV SESD, ATHENS, GA VOLATILES SAMPLE ANALYSIS** Production Date: 06/20/2003 14:10 Produced by: Goddard, Denise Sample 6783 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 14:40 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS14SS / MD No: 1XW3 Org Contractor: LIBRTY D No: 1XW3 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS ANALYTE** UG/KG Dibromochloromethane 11 U UG/KG Dichlorodifluoromethane 11 U 11 U UG/KG Chloromethane 11 U UG/KG 1,2-Dibromoethane (EDB) 11 U UG/KG Vinyl Chloride 11 U UG/KG Chlorobenzene 11 UJ Bromomethane 11 U Ethyl Benzene UG/KG UG/KG 11 U UG/KG Chloroethane 11 U UG/KG Total Xvlenes 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene 11 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 11 U UG/KG Bromoform 11 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG Isopropylbenzene 75 J UG/KG Acetone 1.1.2.2-Tetrachloroethane 11 U UG/KG 11 U UG/KG Carbon Disulfide 1.3-Dichlorobenzene 11 Ú UG/KG 11 U UG/KG Methyl Acetate 11 U UG/KG 1,4-Dichlorobenzene Methylene Chloride 11 U UG/KG 11 U UG/KG 1.2-Dichlorobenzene 11 U trans-1.2-Dichloroethene UG/KG 11 U UG/KG 1.2-Dibromo-3-Chloropropane (DBCP) Methyl T-Butyl Ether (MTBE) 11 U UG/KG 11 U UG/KG 1.2.4-Trichlorobenzene 1.1-Dichloroethane 11 U UG/KG NA UG/KG 1,2,3-Trichlorobenzene cis-1.2-Dichloroethene 11 U UG/KG 14 % Moisture UG/KG Methyl Ethyl Ketone 11 U NA UG/KG Bromochloromethane 11 U UG/KG Chloroform 11 U UG/KG 1.1.1-Trichloroethane

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample **6783** 

6783 FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

i-logodd, ivid

ld/Station: GS14SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

MD No: 1XW3 D No: 1XW3 Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:40

Ending:

RESULTS UNITS

**ANALYTE** 

12 NJ UG/KG ACETALDEHYDE

7 J UG/KG UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

UG/KG

UG/KG

12 U

12 U

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise 6781 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:13 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS02SS / MD No: 1XW1 **Org Contractor: LIBRTY** D No: 1XW1 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS

					אואם	TIEF OFFIED ON DITH WEIGHT BA	.010	
RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE	
12 U	UG/KG	Dichlorodifluoromethane	3		12 U	UG/KG	Dibromochloromethane	•
12 U	UG/KG	Chloromethane	!		12 U	. UG/KG	1,2-Dibromoethane (EDB)	
12 U	UG/KG	Vinyl Chloride			12 U	UG/KG	Chlorobenzene	
12 UJ	UG/KG	Bromomethane			12 U	UG/KG	Ethyl Benzene	e e
12 U	UG/KG	Chloroethane	İ		12 U	UG/KG	Total Xylenes	•
12 U	UG/KG	Trichlorofluoromethane			12 U	UG/KG	Styrene	•
12 U	UG/KG	1,1-Dichloroethene (1,1-			12 U	UG/KG	Bromoform	
12 U	UG/KG	1,1,2-Trichloro-1,2,2-Trif	luoroethane (Freon 113)		-12 U	UG/KG	Isopropylbenzene	
180 J	UG/KG	Acetone	,		12 U	UG/KG	1,1,2,2-Tetrachloroethane	
12 U	UG/KG	Carbon Disulfide			12 U	UG/KG	1,3-Dichlorobenzene	
12 U	UG/KG	Methyl Acetate	1		· 12 U	UG/KG	1,4-Dichlorobenzene	
12 U	UG/KG	Methylene Chloride		•	· 12 U	UG/KG	1,2-Dichlorobenzene	
12 U	UG/KG	trans-1,2-Dichloroethene			12 U	UG/KG	1,2-Dibromo-3-Chloropropane (Di	BCP)
12 U	UG/KG	Methyl T-Butyl Ether (M	ΓBE)		12 U	UG/KG	1,2,4-Trichlorobenzene	
12 U	UG/KG	1,1-Dichloroethane	'		NA	UG/KG	1,2,3-Trichlorobenzene	
12 U	UG/KG	cis-1,2-Dichloroethene		•	23	%	% Moisture	
21	UG/KG	Methyl Ethyl Ketone		i			•	
NA	UG/KG	Bromochloromethane						
12 U	UG/KG	Chloroform						
12 U	UG/KG	1,1,1-Trichloroethane						•
12 U	UG/KG	Cyclohexane			·.			•
12 U	UG/KG	Carbon Tetrachloride			•	• •		
12 U	UG/KG	Benzene						. •
12 U	UG/KG	1,2-Dichloroethane					•	
12 U	UG/KG	Trichloroethene (Trichlo	roethylene)	•				
12 U	UG/KG	Methylcyclohexane						
12 U	UG/KG	1,2-Dichloropropane						
12 U	UG/KG	Bromodichloromethane				•		
12 U	UG/KG	cis-1,3-Dichloropropene			•		•	
12 U	UG/KG	Methyl Isobutyl Ketone						
12 U	UG/KG	Toluene					•	
12 U	UG/KG	trans-1,3-Dichloroproper	ne					•
12 U	UG/KG	1,1,2-Trichloroethane						•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

Tetrachloroethene (Tetrachloroethylene)

Methyl Butyl Ketone

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6781 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:13

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF Id/Station: GS02SS / Case No: 31635

MD No: 1XW1

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

D No: 1XW1

Org Contractor: LIBRTY

**RESULTS UNITS** 51 J UG/KG

**ANALYTE** 

2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

 Sample	<b>6779</b> F	Y <b>2003</b> Project: <b>03-0</b> 4	174					uced by: Goddard, Denise			_
Volatiles	Soon		•		•		Requ	estor:			
							Proje	ct Leader: BSTRIGGO			
		Creosoting Flo	wood, MS				Begin	ning: 04/22/2003 13:40			
Program:			Case No: 31635				Ending:				
ld/Station:	GS15SS /		MD No: 1XT9	Inorg Cont	ractor: S	ENTIN					
Media: SU	RFACE SC	OIL (0" - 12")	D No: 1XT9	Org Contra	actor: LIE	BRTY	DATA	A REPORTED ON DRY WEIGHT B	3ASIS		
 RESULTS	UNITS	ANALYTE			•	RESULTS	UNITS	ANALYTE			_
10 U <sub>.</sub>	UG/KG	Dichlorodifluoromethane	<del> </del>			10 U	UG/KG	Dibromochloromethane			
10 U	UG/KG	Chloromethane				10 U	UG/KG	1,2-Dibromoethane (EDB)			
10 U	UG/KG	Vinyl Chloride		÷		10 U	UG/KG	Chlorobenzene			
10 UJ	UG/KG	Bromomethane			:	10 U	UG/KG	Ethyl Benzene		•	
10 U 10 U	UG/KĞ UG/KG	Chloroethane Trichlorofluoromethane	(Froop 11)			10 U 10 U	UG/KG UG/KG	Total Xylenes			
10 U	UG/KG	1,1-Dichloroethene (1,1-	Dichloroethylone)			10 U	UG/KG	Styrene Bromoform	•	•	
10 U	UG/KG		(luoroethane (Freon 113)	•	•	10 U	UG/KG	Isopropylbenzene			
100 J	UG/KG	Acetone				10 U	UG/KG	1,1,2,2-Tetrachloroethane		•	
10 U	UG/KG	Carbon Disulfide				10 U	UG/KG	1.3-Dichlorobenzene			
3 J	UG/KG	Methyl Acetate				10 U	UG/KG	1,4-Dichlorobenzene			
10 U	UG/KG	Methylene Chloride				10 U	UG/KG	1,2-Dichlorobenzene			
10 U	UG/KG	trans-1,2-Dichloroethen				10 U -	UG/KG	1,2-Dibromo-3-Chloropropane (	DBCP)	•	
10 U	UG/KG	Methyl T-Butyl Ether (M	TBE)			10 U	UG/KG	1,2,4-Trichlorobenzene		•	
10 U	UG/KG	1,1-Dichloroethane					UG/KG	1,2,3-Trichlorobenzene			
10 U 12	UG/KG UG/KG	cis-1,2-Dichloroethene				16	%	% Moisture			
NA		Methyl Ethyl Ketone Bromochloromethane					•			•	
10 U	UG/KG	Chloroform			•						
10 U	UG/KG	1,1,1-Trichloroethane			•						
10 U	UG/KG	Cyclohexane									
10 U	UG/KG	Carbon Tetrachloride						•			
10 U	UG/KG	Benzene									
10 U	UG/KG	1,2-Dichloroethane	•								
10 U	UG/KG	Trichloroethene (Trichlo	roethylene)		•						
.10 U	UG/KG	Methylcyclohexane		•				•			
10 U	UG/KG	1,2-Dichloropropane						•			
10 U	UG/KG	Bromodichloromethane									
10 U 10 U	UG/KG UG/KG	cis-1,3-Dichloropropene				-			•	•	
10 U	UG/KG	Methyl Isobutyl Ketone Toluene									
10 U	UG/KG	trans-1,3-Dichloroproper	ne ,		•						
10 U	UG/KG	1,1,2-Trichloroethane						·			
10 U	UG/KG	Tetrachloroethene (Tetra	achloroethylene)								
10 U	UG/KG	Methyl Butyl Ketone	*					·			
			•	-							

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Program: SF

6779 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:40

Ending:

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

MD No: 1XT9

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

D No: 1XT9

Org Contractor: LIBRTY

**RESULTS UNITS ANALYTE** 

> 24 NJ UG/KG 10 J UG/KG

ld/Station: GS15SS /

**ACETALDEHYDE** UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Project: 03-0474 Sample 6778 FY **2003** Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:55 Program: SF Case No: 31635 Ending: Id/Station: GS23SS / MD No: 1XT8 Inorg Contractor: SENTIN Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT8 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS RESULTS UNITS ANALYTE** ANALYTE 13 U UG/KG Dichlorodifluoromethane 13 U UG/KG Dibromochloromethane 13 U UG/KG Chloromethane 13 U UG/KG 1,2-Dibromoethane (EDB) 13 U UG/KG Vinyl Chloride 13 U UG/KG Chlorobenzene 13 UJ Bromomethane UG/KG 13 U UG/KG Ethyl Benzene 13 U UG/KG Chloroethane 13 U Total Xylenes UG/KG 13 U UG/KG Trichlorofluoromethane (Freon 11) 13 U UG/KG Styrene 13 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 13 U UG/KG Bromoform 13 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 13 U UG/KG Isopropylbenzene 130 J UG/KG Acetone 13 U 1,1,2,2-Tetrachloroethane UG/KG 13 U UG/KG Carbon Disulfide 13 U UG/KG 1.3-Dichlorobenzene 13 U Methyl Acetate 1.4-Dichlorobenzene UG/KG 13 U UG/KG 13 U UG/KG Methylene Chloride 13 U UG/KG 1,2-Dichlorobenzene 13 U trans-1,2-Dichloroethene UG/KG 13 U UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 13 U UG/KG Methyl T-Butyl Ether (MTBE) 13 U 1.2.4-Trichlorobenzene UG/KG 13 U 1,1-Dichloroethane UG/KG NA UG/KG 1.2.3-Trichlorobenzene 13 U UG/KG cis-1,2-Dichloroethene 26 % Moisture 14 UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 13 U UG/KG Chloroform 13 U UG/KG 1,1,1-Trichloroethane 13 U UG/KG Cyclohexane 13 U UG/KG Carbon Tetrachloride 13 U UG/KG Benzene 13 U UG/KG 1.2-Dichloroethane 13 U UG/KG Trichloroethene (Trichloroethylene) 13 U UG/KG Methylcyclohexane 13 U 1,2-Dichloropropane UG/KG 13 U UG/KG Bromodichloromethane 13 U UG/KG cis-1,3-Dichloropropene 13 U UG/KG Methyl Isobutyl Ketone 13 U UG/KG Toluene 13 U UG/KG trans-1,3-Dichloropropene 13 U UG/KG 1,1,2-Trichloroethane 13 U UG/KG Tetrachloroethene (Tetrachloroethylene) 13 U UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6778 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:55

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS23SS /

MD No: 1XT8

D No: 1XT8

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Media: SURFACE SOIL (0" - 12") **RESULTS UNITS** 

**ANALYTE** 

**ACETALDEHYDE** 

14 NJ UG/KG 30 J UG/KG **3 UNKNOWN COMPOUNDS** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

UG/KG

15 UJ UG/KG.

Produced by: Goddard, Denise Project: 03-0474 Sample **6773** FY 2003 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS13SS / MD No: 1XT3 Media: SURFACE SOIL (0" - 12") Org Contractor: LIBRTY D No: 1XT3 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS UNITS	ANALYTE
15 U	UG/KG	Dichlorodifluoromethane	15 U UG/KG	Dibromochloromethane
15 U	UG/KG	Chloromethane	15 U UG/KG	1,2-Dibromoethane (EDB)
15 U	UG/KG	Vinyl Chloride	15 U UG/KG	Chlorobenzene
15 UJ	UG/KG	Bromomethane	15 U UG/KG	Ethyl Benzene
15 U	UG/KG	Chloroethane	15 U UG/KG	Total Xylenes
15 U	UG/KG	Trichlorofluoromethane (Freon 11)	15 U UG/KG	Styrene
15 U	UG/KG	1,1-Dichloroethene (1,1-Dichloroethylene)	15 U UG/KG	Bromoform
15 U	UG/KG	1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	15 U UG/KG	Isopropylbenzene
200 J	UG/KG	Acetone	15 U UG/KG	1,1,2,2-Tetrachloroethane
15 U	UG/KG	Carbon Disulfide	15 U UG/KG	1,3-Dichlorobenzene
15 U	UG/KG	Methyl Acetate	15 U UG/KG	1,4-Dichlorobenzene
15 U	UG/KG	Methylene Chloride	15 U UG/KG	1,2-Dichlorobenzene
15 U	UG/KG	trans-1,2-Dichloroethene	15 UJ UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)
15 U	UG/KG	Methyl T-Butyl Ether (MTBE)	15 U UG/KG	1,2,4-Trichlorobenzene
15 U	UG/KG	1,1-Dichloroethane	NA UG/KG	1,2,3-Trichlorobenzene
15 U	UG/KG	cis-1,2-Dichloroethene	13 %	% Moisture
15 J	UG/KG	Methyl Ethyl Ketone		
NA	UG/KG	Bromochloromethane	•	
15 U	UG/KG	Chloroform	•	
15 U	UG/KG	1,1,1-Trichloroethane		•
15 U	UG/KG	Cyclohexane		•
15 U	UG/KG	Carbon Tetrachloride		
15 U	UG/KG	Benzene	•	
15 U	UG/KG	1,2-Dichloroethane	•	
15 U	UG/KG	Trichloroethene (Trichloroethylene)		•
15 U	UG/KG	Methylcyclohexane		
15 U	UG/KG	1,2-Dichloropropane	•	
15 U	UG/KG	Bromodichloromethane		,
15 U	UG/KG	cis-1,3-Dichloropropene		
15 UJ	UG/KG	Methyl Isobutyl Ketone		
· 15 U	UG/KG	Toluene		
15 U	UG/KG	trans-1,3-Dichloropropene		
15 U	UG/KG	1,1,2-Trichloroethane		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Tetrachloroethene (Tetrachloroethylene)

Methyl Butyl Ketone

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6773 FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS13SS /

MD No: 1XT3

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XT3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:15

Ending:

RESULTS UNITS 16 NJ UG/KG **ANALYTE** 

**ACETALDEHYDE** 

13 J

UG/KG

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

UG/KG

Methyl Butyl Ketone

Produced by: Goddard, Denise 6772 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:00 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS01SS / MD No: 1XT2 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT2 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS RESULTS UNITS ANALYTE ANALYTE** 11 U UG/KG Dichlorodifluoromethane 11 U UG/KG Dibromochloromethane UG/KG Chloromethane 11 U UG/KG 1.2-Dibromoethane (EDB) 11 U 11 U 11 U ÚG/KG Vinvl Chloride UG/KG Chlorobenzene 11 UJ UG/KG Bromomethane 11 U UG/KG Ethyl Benzene 11 U UG/KG Chloroethane 11 U UG/KG Total Xylenes 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene: 11 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 11 U UG/KG Bromoform 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG 11 U UG/KG Isopropylbenzene 100 J UG/KG Acetone 11 U UG/KG 1.1.2.2-Tetrachloroethane 11 U UG/KG Carbon Disulfide 1.3-Dichlorobenzene 11 U UG/KG 11 U UG/KG Methyl Acetate 11 U UG/KG 1.4-Dichlorobenzene 11 U UG/KG Methylene Chloride 11 U UG/KG 1,2-Dichlorobenzene trans-1.2-Dichloroethene 11 U UG/KG 11 U UG/KG 1.2-Dibromo-3-Chloropropane (DBCP) 11 U UG/KG Methyl T-Butyl Ether (MTBE) 11 U UG/KG 1.2.4-Trichlorobenzene 11 U UG/KG 1.1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene cis-1.2-Dichloroethene 11 U UG/KG 14 % Moisture 11 U UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 11 U UG/KG Chloroform 1,1,1-Trichloroethane 11 U UG/KG 11 U UG/KG Cyclohexane UG/KG Carbon Tetrachloride 11 U 11 U UG/KG Benzene 11 U UG/KG 1,2-Dichloroethane Trichloroethene (Trichloroethylene) 11 U UG/KG 11 U UG/KG Methylcyclohexane 11 U UG/KG 1,2-Dichloropropane 11 U Bromodichloromethane UG/KG 11 U cis-1,3-Dichloropropene UG/KG 11 U UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene 11 U UG/KG trans-1,3-Dichloropropene 11 U UG/KG 1,1,2-Trichloroethane Tetrachloroethene (Tetrachloroethylene) 11 U UG/KG

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

**MISCELLANEOUS COMPOUNDS** 

Project: 03-0474

Produced by: Goddard, Denise

6772 FY 2003

Requestor:

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO

Program: SF

Case No: 31635

Beginning: 04/22/2003 09:00

Ending:

Id/Station: GS01SS / Media: SURFACE SOIL (0" - 12") MD No: 1XT2 D No: 1XT2

Inorg Contractor: SENTIN Org Contractor: LIBRTY

**RESULTS UNITS** UG/KG 19 J

ANALYTE

2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6768 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:05

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF Id/Station: GS21SS / Case No: 31635

Media: SURFACE SOIL (0" - 12")

Inorg Contractor: SENTIN MD No: 1XS8 Org Contractor: LIBRTY D No: 1XS8

**RESULTS UNITS ANALYTE** 47 NJ UG/KG

**ACETALDEHYDE PENTANAL** 

7 NJ UG/KG

12 J UG/KG UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Sample 6768 FY **2003** Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 12:05 Program: SF Case No: 31635 Ending: Id/Station: GS21SS / Inorg Contractor: SENTIN MD No: 1XS8 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XS8 DATA REPORTED ON DRY WEIGHT BASIS

	·		1						
RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE	<del></del>	
12 U	UG/KG	Dichlorodifluoromethane			12 U	UG/KG	Dibromochloromethane		
12 U	UG/KG	Chloromethane	i		12 U	UG/KG	1,2-Dibromoethane (EDB	) .	
12 U	UG/KG	Vinyl Chloride			12 U	UG/KG	Chlorobenzene	•	
12 UJ	UG/KG	Bromomethane			12 U	UG/KG	Ethyl Benzene	•	
12 U	UG/KG	Chloroethane	i		· 12 U	UG/KG	Total Xylenes	-	
12 U	UG/KG	Trichlorofluoromethane	(Freon 11)		12 U	UG/KG	Styrene		
12 U	UG/KG	1,1-Dichloroethene (1,1-			12 U	UG/KG	Bromoform		
12 U	UG/KG	1,1,2-Trichloro-1,2,2-Trifl	luoroethane (Freon 113)		12 U	UG/KG	Isopropylbenzene		
270 J	UG/KG	Acetone	i		12 U	UG/KG	1,1,2,2-Tetrachloroethane	و	
12 U	UG/KG	Carbon Disulfide			12 U	UG/KG	1,3-Dichlorobenzene		
12 U	UG/KG	Methyl Acetate	<del>•</del> <u>•</u> •		12 U	UG/KG	1,4-Dichlorobenzene		
12 U	UG/KG	Methylene Chloride	1		12 U	UG/KG	1,2-Dichlorobenzene		
12 U	UG/KG	trans-1,2-Dichloroethene			12 UJ	UG/KG	1,2-Dibromo-3-Chloroprop	oane (DBCP)	
12 U	UG/KG	Methyl T-Butyl Ether (MT			12 U	UG/KG	1,2,4-Trichlorobenzene	, a (220. )	
. 12 U	UG/KG	1,1-Dichloroethane		• .	NA	UG/KG	1,2,3-Trichlorobenzene		
12 U	UG/KG	cis-1,2-Dichloroethene	; ! !		22	%	% Moisture	,	
31 J	UG/KG	Methyl Ethyl Ketone							
NA	UG/KG	Bromochloromethane							
12 U	UG/KG	Chloroform					×	•	
12 U	UG/KG	1,1,1-Trichloroethane							
12 U	UG/KG	Cyclohexane					•		
12 U	UG/KG	Carbon Tetrachloride							•
12 U	UG/KG	Benzene	<u> </u>		•				
12 U	UG/KG	1,2-Dichloroethane	•	•	•				
12 U	UG/KG	Trichloroethene (Trichlor	pethylene)				•		
12 U	UG/KG	Methylcyclohexane		·			·		
12 U	UG/KG	1,2-Dichloropropane			•				
12 U	UG/KG	Bromodichloromethane							
12 U	UG/KG	cis-1,3-Dichloropropene	•				•		
12.UJ	UG/KG	Methyl Isobutyl Ketone		•					•
12 U	UG/KG	Toluene	•						
12 U	UG/KG	trans-1,3-Dichloropropen	e						
12 U	UG/KG	1,1,2-Trichloroethane				•			
12 Ú	UG/KG	Tetrachloroethene (Tetra	chloroethylene)					•	
12 UJ	UG/KG	Methyl Butyl Ketone	,,	•					
•.		, ,							

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6766 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

**MISCELLANEOUS COMPOUNDS** Facility: Gulf States Creosoting

Project Leader: BSTRIGGO

Flowood, MS

Program: SF

16 J

Id/Station: GS16SS /

Case No: 31635

Beginning: 04/22/2003 11:30

Media: SURFACE SOIL (0" - 12")

MD No: 1XS6 D No: 1XS6

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Ending:

Requestor:

**RESULTS UNITS ANALYTE** 31 NJ UG/KG

HEXANAL UG/KG

2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Project: 03-0474 Requestor:

6766 **Volatiles Scan** 

Facility: Gulf States Creosoting

FY 2003

Program: SF

Sample

ld/Station: GS16SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635

MD No: 1XS6 D No: 1XS6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

Project Leader: BSTRIGGO

Beginning: 04/22/2003 11:30

Ending:

RESULTS	UNITS	ANALYTE	RESULTS UNITS	ANALYTE
-12 U	UG/KG	Dichlorodifluoromethane	12 U UG/KG	Dibromochloromethane
12 U	UG/KG	Chloromethane	12 U UG/KG	1,2-Dibromoethane (EDB)
12 U	UG/KG	Vinyl Chloride	12 U UG/KG	Chlorobenzene
12 UJ	UG/KG	Bromomethane	12 U UG/KG	Ethyl Benzene
12 U	UG/KG	Chloroethane	12 U UG/KG	Total Xylenes
12 U	UG/KG	Trichlorofluoromethane (Freon 11)	12 U UG/KG	Styrene
12 U	UG/KG	1,1-Dichloroethene (1,1-Dichloroethylene)	12 U UG/KG	Bromoform
12 U	UG/KG	1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113)	12 U UG/KG	Isopropylbenzene
56 J	UG/KG	Acetone	12 U UG/KG	1,1,2,2-Tetrachloroethane
12 U	UG/KG	Carbon Disulfide	12 U UG/KG	1,3-Dichlorobenzene
12 U	UG/KG	Methyl Acetate	12 U UG/KG	1,4-Dichlorobenzene
12 U	UG/KG	Methylene Chloride	12 U UG/KG	1,2-Dichlorobenzene
. 12 U	UG/KG	trans-1,2-Dichloroethene	12 U UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)
12 U	UG/KG	Methyl T-Butyl Ether (MTBE)	12 U UG/KG	1,2,4-Trichlorobenzene
12 U	UG/KG	1,1-Dichloroethane	NA UG/KG	1,2,3-Trichlorobenzene
12 U	UG/KG	cis-1,2-Dichloroethene	21 %	% Moisture
12 U	UG/KG	Methyl Ethyl Ketone	•	
NA	UG/KG	Bromochloromethane	• •	•
12 U	UG/KG	Chloroform		·
12 U	UG/KG	1,1,1-Trichloroethane		
12 U	UG/KG	Cyclohexane		
12 U	UG/KG	Carbon Tetrachloride		
	UG/KG	Benzene		
. — –	UG/KG	1,2-Dichloroethane		
12 U	UG/KG	Trichloroethene (Trichloroethylene)		
	UG/KG UG/KG	Methylcyclohexane		
	UG/KG	1,2-Dichloropropane		
	UG/KG	Bromodichloromethane	•	
	UG/KG	cis-1,3-Dichloropropene Methyl Isobutyl Ketone	•	•
	UG/KG	Toluene		
	UG/KG	trans-1,3-Dichloropropene		•
	UG/KG	1,1,2-Trichloroethane	•	
	UG/KG	Tetrachloroethene (Tetrachloroethylene)		
	UG/KG	Methyl Butyl Ketone		,
12.0	Junu	Metriyi Dutyi Netorie		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyze analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Ending:

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Project: 03-0474 6762 FY 2003 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 10:45

Program: SF Case No: 31635 Inorg Contractor: SENTIN Id/Station: GS20SS / MD No: 1XS2

Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XS2 DATA REPORTED ON DRY WEIGHT BASIS

			,	1		 		5,,,,,	THE OTHER OTTE	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,			
	RESULTS  12 U  12 U  12 U  12 U  12 U  12 U  12 U  12 U  12 U  12 U  12 U  12 U  12 U  13 U	UNITS UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	ANALYTE Dichlorodifluoromethane Chloromethane Vinyl Chloride Bromomethane Chloroethane Trichlorofluoromethane 1,1-Dichloroethene (1,1- 1,1,2-Trichloro-1,2,2-Trif Acetone Carbon Disulfide	(Freon 11) Dichloroethylene)			RESULTS 12 U 12 U 12 U 12 U 12 U 12 U 12 U 12 U	UNITS UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	ANALYTE Dibromochlorome 1,2-Dibromoetha Chlorobenzene Ethyl Benzene Total Xylenes Styrene Bromoform isopropylbenzene 1,1,2,2-Tetrachlo 1,3-Dichlorobenz	ne (EDB)			
	12 U	UG/KG	Methyl Acetate				12 U 12 U	UG/KG UG/KG	1,3-Dichlorobenz				•
·	12 U 12 U 12 U 12 U 12 U 12 U	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG	Methylene Chloride trans-1,2-Dichloroethene Methyl T-Butyl Ether (MT 1,1-Dichloroethane cis-1,2-Dichloroethene Methyl Ethyl Ketone				12 U 12 U 12 U 12 U NA 21	UG/KG UG/KG UG/KG UG/KG	1,2-Dichlorobenz 1,2-Dibromo-3-Cl 1,2,4-Trichlorobe 1,2,3-Trichlorobe % Moisture	ene nioropropa nzene	ne (DBCP)		
	NA 12 U	UG/KG UG/KG	Bromochloromethane Chloroform										
	12 U	UG/KG	1,1,1-Trichloroethane					•					
	12 U	UG/KG	Cyclohexane										
	12 U 12 U	UG/KG UG/KG	Carbon Tetrachloride Benzene									•	
	12 U	UG/KG	1,2-Dichloroethane		•				•				•
	12 U	UG/KG	Trichloroethene (Trichlor	pethylene)									
	12 U 12 U	UG/KG UG/KG	Methylcyclohexane 1,2-Dichloropropane	·									
	12 U	UG/KG	Bromodichloromethane									. '	
	12 U	UG/KG	cis-1,3-Dichloropropene			•							
	12 U	UG/KG UG/KG	Methyl Isobutyl Ketone Toluene								-		
	12 U 12 U	UG/KG	trans-1,3-Dichloropropen	i e							•		
	12 U	UG/KG	1,1,2-Trichloroethane	i i									
	12 U	UG/KG	Tetrachloroethene (Tetra	chloroethylene)			٠.						
	12 U	UG/KG	Methyl Butyl Ketone										

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6760 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

12 J

ld/Station: GS17SS /

MD No: 1XS0

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XS0

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:05

Ending:

**RESULTS UNITS ANALYTE** 

UG/KG

38 NJ UG/KG 7 NJ UG/KG **ACETALDEHYDE** PROPANAL, 2-METHYL-

UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6760 FY 2003 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 10:05 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS17SS /-MD No: 1XS0 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XS0 DATA REPORTED ON DRY WEIGHT BASIS

				<del>'</del>				<u></u>		
R	ESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE		•
	13 U	UG/KG	Dichlorodifluoromethane			13 U	UG/KG	Dibromochloromethane		
	13 U	UG/KG	Chloromethane		•	13 U	UG/KG	1,2-Dibromoethane (EDB	)	
	13 U	UG/KG	Vinyl Chloride			13 U	UG/KG	Chlorobenzene	·	
	13 UJ	UG/KG	Bromomethane			13 U	UG/KG	Ethyl Benzene		
	13 U	UG/KG	Chloroethane			13 U	UG/KG	Total Xylenes		
	13 U	UG/KG	Trichlorofluoromethane (Free	n 11)		13 U	UG/KG	Styrene		
	13 U	UG/KG	1,1-Dichloroethene (1,1-Dichle		•	13 U	UG/KG	Bromoform		·
	13 U	UG/KG	1,1,2-Trichloro-1,2,2-Trifluoro		•	13 U	UG/KG	Isopropylbenzene		
	260 J	UG/KG	Acetone			13 Ü	UG/KG	1,1,2,2-Tetrachloroethane	4	•
	13 U	UG/KG	Carbon Disulfide			13 U	UG/KG	1,3-Dichlorobenzene	•	
	13 U	UG/KG	Methyl Acetate			13 U	UG/KG	1,4-Dichlorobenzene		
	13 U	UG/KG	Methylene Chloride		•	13 U	·UG/KG	1,2-Dichlorobenzene		
	13 U	UG/KG	trans-1,2-Dichloroethene			13 UJ	UG/KG	1,2-Dibromo-3-Chloroprop	oane (DBCP)	
	13 U	UG/KG	Methyl T-Butyl Ether (MTBE)			13 U	UG/KG	1,2,4-Trichlorobenzene		
	13 U	UG/KG	1,1-Dichloroethane	•		NA	UG/KG	1,2,3-Trichlorobenzene		
	13 U	UG/KG	cis-1,2-Dichloroethene			21	%	% Moisture		
	28 J	UG/KG	Methyl Ethyl Ketone					70 1112121		
	NA	UG/KG	Bromochloromethane							
	13 U	UG/KG	Chloroform							
	13 U	UG/KG	1,1,1-Trichloroethane		•		•			
	13 U	UG/KG	Cyclohexane					•		
	13 U	UG/KG	Carbon Tetrachloride							
	13 U	UG/KG	Benzene							
	13 U	UG/KG	1,2-Dichloroethane							
	13 U	UG/KG	Trichloroethene (Trichloroethy	lene)	•			. <del>-</del>		
•	13 U	UG/KG	Methylcyclohexane	,					•	
	13 U	UG/KG	1,2-Dichloropropane				•			
	13 U	UG/KG	Bromodichloromethane			•	·			
	13 U	UG/KG	cis-1,3-Dichloropropene			•	•			•
	13 UJ	UG/KG	Methyl Isobutyl Ketone							
	13 U	UG/KG	Toluene							
	13 Ü	UG/KG	trans-1,3-Dichloropropene							
	13 U	UG/KG	1,1,2-Trichloroethane							
	13 U	UG/KG	Tetrachloroethene (Tetrachlor	oethvlene)				•		
	13 UJ	UG/KG	Methyl Butyl Ketone				4.			
					•					
								•		

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6758 FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS19SS /

Media: SURFACE SOIL (0" - 12")

MD No: 1XR8

D No: 1XR8

Case No: 31635

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:30

Endina:

**RESULTS UNITS** 18 NJ UG/KG

**ANALYTE** 

**ACETALDEHYDE** 

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Sample 6758 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:30 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS19SS / MD No: 1XR8 Org Contractor: LIBRTY D No: 1XR8 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS RESULTS UNITS** ANALYTE **ANALYTE** UG/KG 11 U UG/KG Dichlorodifluoromethane 11 U Dibromochloromethane 1,2-Dibromoethane (EDB) 11 U UG/KG Chloromethane 11 U UG/KG 11 U UG/KG Vinyl Chloride 11 U UG/KG Chlorobenzene 11 UJ UG/KG Bromomethane 11 U UG/KG Ethyl Benzene UG/KG Chloroethane Total Xylenes 11 U 11 U UG/KG UG/KG Trichlorofluoromethane (Freon 11) Styrene 11 U 11 U UG/KG 1.1-Dichloroethene (1.1-Dichloroethylene) UG/KG Bromoform 11 U 11 U UG/KG 11 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG Isopropylbenzene 180 J

UG/KG Acetone 11 U UG/KG 1,1,2,2-Tetrachloroethane 11 U UG/KG Carbon Disulfide 11 U UG/KG 1,3-Dichlorobenzene Methyl Acetate 3 J UG/KG 11 U UG/KG 1.4-Dichlorobenzene Methylene Chloride 11 U UG/KG 11 U UG/KG 1,2-Dichlorobenzene trans-1.2-Dichloroethene 11 U UG/KG 11.UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) UG/KG Methyl T-Butyl Ether (MTBE) 11 U 11 U UG/KG 1,2,4-Trichlorobenzene 11 U UG/KG 1.1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene cis-1,2-Dichloroethene -11 U UG/KG % Moisture Methyl Ethyl Ketone 17 J UG/KG NA UG/KG Bromochloromethane 11 U UG/KG Chloroform

1,2-Dichloroethane UG/KG Trichloroethene (Trichloroethylene) 11 U 11 U UG/KG Methylcyclohexane 11 U UG/KG 1,2-Dichloropropane 11 U UG/KG Bromodichloromethane

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

11 U

11 U

11 U

11 U

11 U

11 U UG/KG cis-1,3-Dichloropropene 11 UJ UG/KG Methyl Isobutyl Ketone UG/KG Toluene

11 U UG/KG trans-1,3-Dichloropropene 11 U 11 U UG/KG 1.1.2-Trichloroethane

11 U UG/KG Tetrachloroethene (Tetrachloroethylene)

1,1,1-Trichloroethane

Carbon Tetrachloride

Cyclohexane

Benzene

11 UJ UG/KG Methyl Butyl Ketone

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6756 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS18SS /

Media: SURFACE SOIL (0" - 12")

MD No: 1XR6

D No: 1XR6

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:40

Ending:

RESULTS UNITS **ANALYTE** 

8 J UG/KG UNKNOWN COMPOUND

20 NJ UG/KG **HEXANAL** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

13 U

13 U

13 U

13 U

13 U

13 U

13 U

13 U

13 U

13 U

13 U

13 UJ

13 UJ

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

Production Date: 06/20/2003 14:10 Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6756 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 08:40 Program: SF Case No: 31635 Ending: Id/Station: GS18SS / Inorg Contractor: SENTIN MD No: 1XR6 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XR6 DATA REPORTED ON DRY WEIGHT BASIS **ANALYTE RESULTS UNITS RESULTS UNITS** ANALYTE 13 U UG/KG Dichlorodifluoromethane 13 U UG/KG Dibromochloromethane 13 U Chloromethane 13 U UG/KG UG/KG 1.2-Dibromoethane (EDB) Vinyl Chloride 13 U UG/KG 13 U UG/KG Chlorobenzene 13 UJ UG/KG Bromomethane 13 U UG/KG Ethyl Benzene 13 U UG/KG Chloroethane 13 U UG/KG Total Xvlenes 13 U UG/KG Trichlorofluoromethane (Freon 11) 13 U UG/KG Styrene 13 U -UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 13 U UG/KG Bromoform UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 13 U 13 U UG/KG

130 J UG/KG Acetone 13 U UG/KG Carbon Disulfide 13 U UG/KG Methyl Acetate 13 U Methylene Chloride UG/KG 13 U UG/KG trans-1.2-Dichloroethene 13 U UG/KG Methyl T-Butyl Ether (MTBE) 13 U UG/KG 1.1-Dichloroethane cis-1,2-Dichloroethene 13 U UG/KG Methyl Ethyl Ketone 14 J UG/KG NA UG/KG Bromochloromethane 13 U UG/KG Chloroform 13 U UG/KG 1.1.1-Trichloroethane UG/KG 13 U Cyclohexane

Carbon Tetrachloride

Trichloroethene (Trichloroethylene)

Tetrachloroethene (Tetrachloroethylene)

1,2-Dichloroethane

Methylcyclohexane

1,2-Dichloropropane

Bromodichloromethane

cis-1,3-Dichloropropene

Methyl Isobutyl Ketone

1,1,2-Trichloroethane

Methyl Butyl Ketone

trans-1,3-Dichloropropene

Benzene

Toluene

Isopropylbenzene 13 U UG/KG 1,1,2,2-Tetrachloroethane 13 U UG/KG 1,3-Dichlorobenzene 13 U UG/KG 1.4-Dichlorobenzene 13 U UG/KG 1.2-Dichlorobenzene 13 UJ UG/KG 1.2-Dibromo-3-Chloropropane (DBCP) 13 U UG/KG 1,2,4-Trichlorobenzene NA UG/KG 1,2,3-Trichlorobenzene 23 % % Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

12 U

12 U

12 U 12 U

12 U

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG 12 UJ UG/KG

12 UJ UG/KG

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise 6803 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 17:54 Program: SF Case No: 31635 Endina: Id/Station: GS01TB / Org Contractor: LIBRTY Media: TRIP BLANK - SOIL D No: 1XX8 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS ANALYTE** Dichlorodifluoromethane 12 U UG/KG Dibromochloromethane 12 U UG/KG 12 U UG/KG 1.2-Dibromoethane (EDB) 12 U UG/KG Chloromethane -12 U UG/KG Vinvl Chloride 12 U UG/KG Chlorobenzene 12 U UG/KG Ethyl Benzene 12 U UG/KG Bromomethane 12 U - 12 U UG/KG Chloroethane UG/KG Total Xylenes 12 U UG/KG Trichlorofluoromethane (Freon 11) 12 U UG/KG Styrene 1,1-Dichloroethene (1,1-Dichloroethylene) 12 U UG/KG 12 U UG/KG Bromoform 12 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 12 U .UG/KG Isopropylbenzene 1,1,2,2-Tetrachloroethane 19 J UG/KG Acetone 12 U UG/KG 12 U UG/KG Carbon Disulfide 12 U UG/KG 1.3-Dichlorobenzene 12 U UG/KG Methyl Acetate 12 U UG/KG 1.4-Dichloropenzene UG/KG Methylene Chloride 12 U UG/KG 1.2-Dichlorobenzene 12 U UG/KG trans-1,2-Dichloroethene 12 UJ 1,2-Dibromo-3-Chloropropane (DBCP) 12 U UG/KG UG/KG Methyl T-Butyl Ether (MTBE) 12 U 12 U UG/KG 1,2,4-Trichlorobenzene 12 U UG/KG 1,1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene 12 U UG/KG cis-1,2-Dichloroethene % Moisture 12 UJ UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane UG/KG 12 U Chloroform · UG/KG 1.1,1-Trichloroethane 12 U Cyclohexane -12 U .UG/KG 12 U . UG/KG Carbon Tetrachloride 12 U UG/KG Benzene 12 U UG/KG 1,2-Dichloroethane 12 U UG/KG Trichloroethene (Trichloroethylene) UG/KG Methylcyclohexane 12 U 1,2-Dichloropropane 12 U UG/KG

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

Bromodichloromethane

cis-1,3-Dichloropropene

trans-1,3-Dichloropropene

Tetrachloroethene (Tetrachloroethylene)

Methyl Isobutyl Ketone

1.1.2-Trichloroethane

Methyl Butyl Ketone

Toluene

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample 6814 FY 2003 Project: 03-0474

**Volatiles Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

19 UJ

UG/KG

Methyl Butyl Ketone

Case No: 31635

Id/Station: GS02TS /

Media: TRIP BLANK - SOIL

D No: 1XZ2

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/23/2003 15:57

Ending:

Requestor:

DATA REPORTED ON DRY WEIGHT BASIS

wicdia. 11,11	II DEAMY	JOIL		Org Comme	ZOTOT: ETE		DATA REPORTED ON DRY WEIGHT BASIS			
RESULTS	UNITS	ANALYTE	:			RESULTS	UNITS	ANALYTE		
19 U	UG/KG	Dichlorodifluoromethane				19 U	UG/KG	Dibromochloromethane		
19 U	UG/KG	Chloromethane	:			19 U	UG/KG	1,2-Dibromoethane (EDB)	)	
19 U	UG/KG	Vinyl Chloride				19 U	UG/KG	Chlorobenzene		
19 UJ	UG/KG	Bromomethane	1			19 U	UG/KG	Ethyl Benzene	•	
19 U	UG/KG	Chloroethane				19 U	UG/KG	Total Xylenes		
19 U	UG/KG	Trichlorofluoromethane	(Freon 11)	·.		19 U	UG/KG	Styrene		•
19 U	UG/KG	1,1-Dichloroethene (1,1-	Dichloroethylene)			19 U	UG/KG	Bromoform		
19 U	. UG/KG	1,1,2-Trichloro-1,2,2-Trif	luoroethane (Freon 113)		-	19 U	UG/KG	Isopropylbenzene		•
19 UJ	UG/KG	Acetone	:			19 U	UG/KG	1,1,2,2-Tetrachloroethane	<del>)</del>	
- 19 U	UG/KG	Carbon Disulfide	:			19 U	UG/KG	1,3-Dichlorobenzene		
19 U	UG/KG	Methyl Acetate				. 19 U	UG/KG	1,4-Dichlorobenzene		
19 U	UG/KG	Methylene Chloride	:			19 U	UG/KG	1,2-Dichlorobenzene		
19 U	UG/KG	trans-1,2-Dichloroethene				19 UJ	UG/KG	1,2-Dibromo-3-Chloroprop	oane (DBCP)	
19 UJ	UG/KG	Methyl T-Butyl Ether (MT	BE)			19 U	UG/KG	1,2,4-Trichlorobenzene	, ,	·
19 U	UG/KG	1,1-Dichloroethane	ì i			NA	UG/KG	1,2,3-Trichlorobenzene		
19 U	UG/KG	cis-1,2-Dichloroethene		•		0	%	% Moisture		
19 UJ	UG/KG	Methyl Ethyl Ketone	•					• • •		
· NA	UG/KG	Bromochloromethane							•	•
19 U	UG/KG	Chloroform								
19 U	UG/KG	1,1,1-Trichloroethane	-							
19 U	UG/KG	Cyclohexane	:				•			
19 U	UG/KG	Carbon Tetrachloride						•	•	
19 U	UG/KG	Benzene								
19 U	UG/KG	1,2-Dichloroethane								
19 U	UG/KG	Trichloroethene (Trichlor	oethylene)							
19 U	UG/KG	Methylcyclohexane				•				
19 U	UG/KG	1,2-Dichloropropane								
19 U	UG/KG	Bromodichloromethane								
19 U	·UG/KG	cis-1,3-Dichloropropene								
19 UJ	UG/KG	Methyl Isobutyl Ketone	•							
19 U	UG/KG	Toluene							•	
· 19 U	UG/KG	trans-1,3-Dichloropropen	e .	÷						•
·19 U	UG/KG	1,1,2-Trichloroethane	i i							
19 U	UG/KG	Tetrachloroethene (Tetra	chloroethylene)					•		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences, | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6771 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 09:15 Program: SF Case No: 31635 Ending: Id/Station: GS01SB / Inorg Contractor: SENTIN MD No: 1XT1 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XT1 DATA REPORTED ON DRY WEIGHT BASIS RESULTS UNITS **ANALYTE RESULTS UNITS ANALYTE** Dichlorodifluoromethane 11 U UG/KG 11 U UG/KG Dibromochloromethane Chloromethane 11 U UG/KG 11 U UG/KG 1,2-Dibromoethane (EDB) 11 U UG/KG Vinyl Chloride 11 U UG/KG Chlorobenzene 11 UJ UG/KG Bromomethane 11 U UG/KG Ethyl Benzene Chloroethane 11 U UG/KG 11 U UG/KG Total Xylenes 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene UG/KG 1.1-Dichloroethene (1.1-Dichloroethylene) 11 U UG/KG 11 U Bromoform 11 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG Isopropylbenzene 39 UJ UG/KG Acetone 11 U UG/KG 1.1.2.2-Tetrachloroethane 11 U UG/KG Carbon Disulfide 11 U UG/KG 1.3-Dichlorobenzene 11 U UG/KG Methyl Acetate 11 U UG/KG 1,4-Dichlorobenzene 11 U UG/KG Methylene Chloride 11 U UG/KG 1.2-Dichlorobenzene 11 U UG/KG trans-1,2-Dichloroethene 11 UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) · UG/KG 11 U Methyl T-Butyl Ether (MTBE) 11 U UG/KG 1,2,4-Trichlorobenzene 11 U UG/KG 1,1-Dichloroethane NA UG/KG 1,2,3-Trichlorobenzene 11 U UG/KG cis-1,2-Dichloroethene % % Moisture 18 11 UJ UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane UG/KG Chloroform 11 U 11 U UG/KG 1,1,1-Trichloroethane 11 U UG/KG Cyclohexane 11 U UG/KG Carbon Tetrachloride 11 U UG/KG Benzene 11 U UG/KG 1,2-Dichloroethane 11 U UG/KG Trichloroethene (Trichloroethylene) 11 U UG/KG Methylcyclohexane 11 U UG/KG 1.2-Dichloropropane Bromodichloromethane 11 U UG/KG 11 U UG/KG cis-1,3-Dichloropropene 11 UJ UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene 11 U UG/KG trans-1.3-Dichloropropene 11 U UG/KG 1,1,2-Trichloroethane 11 U UG/KG Tetrachloroethene (Tetrachloroethylene) 11 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Produced by: Goddard, Denise 6782 Project: 03-0474 Sample FY 2003 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:34 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS02SB / MD No: 1XW2 Media: SUBSUBFACE SOIL (> 12") D No: 1XW2 Org Contractor: LIBRTY DATA DEDODTED ON DRY WEIGHT DAGIS

Media: SUE	SUHFAC	JE SOIL (> 12")	D No: 1XW2	Org Contractor: Li	BRIT	DATA	REPORTED ON DRY WEIGHT BAS	IS
RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE	
11 U	UG/KG	Dichlorodifluoromethane		•	11 U	UG/KG	Dibromochloromethane	
11 U	UG/KG	Chloromethane			11 U	UG/KG	1,2-Dibromoethane (EDB)	
11 U	UG/KG	Vinyl Chloride	1		11 U	UG/KG	Chlorobenzene	
11 UJ	UG/KG	Bromomethane	· .		11 U	UG/KG	Ethyl Benzene	
11 U	UG/KG	Chloroethane	· ·		11 U	UG/KG	Total Xylenes	
11 U	UG/KG	Trichlorofluoromethane	(Freon 11)		11 U	UG/KG	Styrene	
11 U	UG/KG	1,1-Dichloroethene (1,1-	Dichloroethylene)		11 U	UG/KG	Bromoform	
-11 U	UG/KG	1,1,2-Trichloro-1,2,2-Trif	luoroethane (Freon 113	•	11 U	UG/KG	Isopropylbenzene	•
12 J	UG/KG	Acetone			11 U	UG/KG	1,1,2,2-Tetrachloroethane	
11 U	UG/KG	Carbon Disulfide	·	•	11 U	UG/KG	1,3-Dichlorobenzene	•
. 11 U	UG/KG	Methyl Acetate			11 U	UG/KG	1,4-Dichlorobenzene	
11 U	UG/KG	Methylene Chloride			11 U	UG/KG	1,2-Dichlorobenzene	
11 U	UG/KG	trans-1,2-Dichloroethene			11 U	UG/KG	1,2-Dibromo-3-Chloropropane (DB)	CP)
11 U	UG/KG	Methyl T-Butyl Ether (MT	BE)		11 U	UG/KG	1,2,4-Trichlorobenzene	
	UG/KG	1,1-Dichloroethane			NA	UG/KG	1,2,3-Trichlorobenzene	
·11 U	UG/KG	cis-1,2-Dichloroethene			19	% .	% Moisture	• .
11 U	UG/KG	Methyl Ethyl Ketone		•				
NA	UG/KG	Bromochloromethane		•				
11 U 🐠	UG/KG	Chloroform	•	•				
11 U	UG/KG	1,1,1-Trichloroethane						• •
	UG/KG	Cyclohexane	2					
	UG/KG	Carbon Tetrachloride					•	
	UG/KG	Benzene	•					•
	UG/KG	1,2-Dichloroethane						
	UG/KG	Trichloroethene (Trichlor	oethylene)	•				
	UG/KG	Methylcyclohexane				•		
	UG/KG	1,2-Dichloropropane					•	
	UG/KG	Bromodichloromethane			•			
	UG/KG	cis-1,3-Dichloropropene		•			·	
	UG/KG	Methyl Isobutyl Ketone					· ·	
	UG/KG	Toluene						•
· · · · ·	UG/KG	trans-1,3-Dichloroproper	e				•	•
_	UG/KG	1,1,2-Trichloroethane			•			,
	UG/KG	Tetrachloroethene (Tetra	chloroethylene)					
11 U	UG/KG	Methyl Butyl Ketone	•					•
•			•					

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate, Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

6801 FY 2003 Project: 03-0474 Sample

**Volatiles Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS03SB / Media: SUBSURFACE SOIL (> 12")

MD No: 1XY2 D No: 1XY2

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Beginning: 04/23/2003 07:55 Ending:

Requestor:

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	1		RESULTS	UNITS	ANALYTE	
23 U	UG/KG	Dichlorodifluoromethane			23 U	UG/KG	Dibromochloromethane	
23 U	UG/KG	Chloromethane	٠.		23 U	UG/KG	1,2-Dibromoethane (EDB)	
23 U	UG/KG	Vinyl Chloride	1 :		23 U	UG/KG	Chlorobenzene	
23 UJ	UG/KG	Bromomethane			23 Ŭ	UG/KG	Ethyl Benzene	
23 U	UG/KG	Chloroethane	. :		23 U	UG/KG	Total Xylenes	•
23 U	UG/KG	Trichlorofluoromethane	(Freen 11)		23 U	UG/KG	Styrene	
23 U	UG/KG	1,1-Dichloroethene (1,1-			23 U	UG/KG	Bromoform	
23 U	UG/KG		luoroethane (Freon 113)		23 U	UG/KG	Isopropylbenzene	
35 J	UG/KG	Acetone	doroemane (ricen, rie)		23 Ŭ	UG/KG	1,1,2,2-Tetrachloroethane	
	- UG/KG	Carbon Disulfide	•	•	23 U	UG/KG	1,3-Dichlorobenzene	
23 U	UG/KG	Methyl Acetate			23 U	UG/KG	1,4-Dichlorobenzene	
23 U	UG/KG	Methylene Chloride		•	23 Ú	UG/KG	1,2-Dichlorobenzene	
23 Ú	UG/KG	trans-1,2-Dichloroethene	) 2		23 UJ	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)	
23 UJ	UG/KG	Methyl T-Butyl Ether (M)			23 U	UG/KG	1,2,4-Trichlorobenzene	•
23 U	UG/KG	1,1-Dichloroethane	,		NA NA	UG/KG	1,2,3-Trichlorobenzene	
23 U	UG/KG	cis-1,2-Dichloroethene	•		28	%	% Moisture	
23 UJ	UG/KG	Methyl Ethyl Ketone					, , , , , , , , , , , , , , , , , , , ,	
NA	UG/KG	Bromochloromethane	•	•				
23 U	UG/KG	Chloroform				•	· ·	
23 U	UG/KG	1,1,1-Trichloroethane						
23 U	UG/KG	Cyclohexane						
23 U	UG/KG	Carbon Tetrachloride					•.	
.23 U	UG/KG	Benzene					•	
23 U	UG/KG	1,2-Dichtoroethane		•			•	
23 U	UG/KG	Trichloroethene (Trichlor	oethylene)					
23 U	UG/KG	Methylcyclohexane	-					
23 U	UG/KG	1,2-Dichloropropane						
23 U	UG/KG	Bromodichloromethane		5				
23 U	UG/KG	cis-1,3-Dichloropropene						
23 UJ	UG/KG	Methyl Isobutyl Ketone		i				
23 U	UG/KG	Toluene						
23 U	UG/KG	trans-1,3-Dichloroproper	ie					
23 U	UG/KG	1,1,2-Trichloroethane						
23 U	UG/KG	Tetrachloroethene (Tetra	chloroethylene)			•		
23 UJ	UG/KG	Methyl Butyl Ketone						

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample 6808 FY 2003 Project: 03-0474 **Volatiles Scan** 

Facility: Gulf States Creosoting

Program: SF Id/Station: GS04SB /

Media: SUBSURFACE SOIL (> 12")

Flowood, MS

Case No: 31635

MD No: 1XY8 Inorg Contractor: SENTIN

Org Contractor: LIBRTY D No: 1XY8

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

				<del></del>	<del></del>				<u> </u>	· · · · · · · · · · · · · · · · · · ·
RESULTS		ANAĻYTĒ				RESULTS		ANALYTE		
15 U	UG/KG	Dichlorodifluoromethane	;			15 U	UG/KG	Dibromochloromethane	•	
15 U	UG/KG	Chloromethane	I			15 U	UG/KG	1,2-Dibromoethane (EDB)		
15 U	UG/KG	Vinyl Chloride	٠.	-		15 U	UG/KG	Chlorobenzene		
15 UJ	UG/KG	Bromomethane	ı			15 U	UG/KG	Ethyl Benzene		
15 U	UG/KG	Chloroethane	i			15 U	UG/KG	Total Xylenes		•
15 U	UG/KG	Trichlorofluoromethane				15 U	UG/KG	Styrene		
15 U	UG/KG	1,1-Dichloroethene (1,1-		•		15 U	UG/KG	Bromoform		
:15 U	UG/KG	1,1,2-Trichloro-1,2,2-Trif	luoroethane (Freon 113)			15 U	UG/KG	Isopropylbenzene		
15 UJ	UG/KG	Acetone				15 U	UG/KG	1,1,2,2-Tetrachloroethane		
15 U	UG/KG	Carbon Disulfide				15 U	UG/KG	1,3-Dichlorobenzene		
15 U	UG/KG	Methyl Acetate				15 U	UG/KG	1,4-Dichlorobenzene		
15 U	UG/KG	Methylene Chloride				15 U	UG/KG	1,2-Dichlorobenzene		
15 U	UG/KG	trans-1,2-Dichloroethene	<b>)</b>	ŧ		15 UJ	UG/KG	1,2-Dibromo-3-Chloropropa	ane (DBCP)	
15 UJ	UG/KG	Methyl T-Butyl Ether (MT	BE)			15 U	UG/KG	1,2,4-Trichlorobenzene	, ,	•
15 U	UG/KG	1,1-Dichloroethane				NA	UG/KG	1,2,3-Trichlorobenzene		
15 U	UG/KG	cis-1,2-Dichloroethene				23	%	% Moisture		
15 UJ	UG/KG	Methyl Ethyl Ketone						•		
NA	UG/KG	Bromochloromethane						•	•	•
15 U	UG/KG	Chloroform								
15 U	UG/KG	1,1,1-Trichloroethane								* .
15 U	UG/KG	Cyclohexane								•
15 U	UG/KG	Carbon Tetrachloride								•
15 U	UG/KG	Benzene						· .	. •	
15 U.	UG/KG	1,2-Dichloroethane				•				
15 U	UG/KG	Trichloroethene (Trichlor	oethylene)							
15 Ü	UG/KG	Methylcyclohexane	<i>,</i>							
15 U	UG/KG	1,2-Dichloropropane						•		•
15 U .	UG/KG	Bromodichloromethane	•				•	•		
15 U	UG/KG	cis-1,3-Dichloropropene								
15 UJ	UG/KG	Methyl Isobutyl Ketone	• .							
15 U	UG/KG	Toluene						•	•	•
15 U	UG/KG	trans-1,3-Dichloropropen	ie							
15 U	UG/KG	1,1,2-Trichloroethane	-							
15 U	UG/KG	Tetrachloroethene (Tetra	chloroethylene)							
15 UJ	UG/KG	Methyl Butyl Ketone								
		,,,,				•		•		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates...

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample -6810 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 09:25 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS05SB / MD No: 1XZ0 Media: SUBSURFACE SOIL (> 12") D No: 1XZ0 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS RESULTS UNITS ANALYTE ANALYTE** . 12 U UG/KG Dichlorodifluoromethane 12 U UG/KG Dibromochloromethane 12 U UG/KG 1,2-Dibromoethane (EDB) 12 U UG/KG Chloromethane UG/KG Vinyl Chloride 12 U UG/KG Chlorobenzene 12 U 12 UJ UG/KG Bromomethane 12 U UG/KG Ethyl Benzene 12 U UG/KG Chloroethane 12 U UG/KG **Total Xylenes** UG/KG Trichlorofluoromethane (Freon 11) 12 U UG/KG Styrene 12 U 1,1-Dichloroethene (1,1-Dichloroethylene) 12 U Bromoform 12 U UG/KG UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) UG/KG 12 U UG/KG Isopropylbenzene 12 U 1.1.2.2-Tetrachloroethane 12 UJ UG/KG Acetone 12 U UG/KG UG/KG Carbon Disulfide 12 U 12 U UG/KG 1.3-Dichlorobenzene 12 U UG/KG Methyl Acetate 12 U UG/KG 1,4-Dichlorobenzene 12 U UG/KG Methylene Chloride 12 U UG/KG 1.2-Dichlorobenzene 12 U UG/KG trans-1,2-Dichloroethene 12 UJ UG/KG 1.2-Dibromo-3-Chloropropane (DBCP) UG/KG Methyl T-Butyl Ether (MTBE) 1.2.4-Trichlorobenzene 12 UJ 12 U UG/KG NA UG/KG 12 U UG/KG 1.1-Dichloroethane 1,2,3-Trichlorobenzene 12 U UG/KG cis-1,2-Dichloroethene 19 % Moisture Methyl Ethyl Ketone 12 UJ UG/KG NA UG/KG Bromochloromethane 12 U UG/KG Chloroform 12 U UG/KG 1,1,1-Trichloroethane 12 U UG/KG Cyclohexane UG/KG Carbon Tetrachloride 12 U 12 U UG/KG Benzene 12 U UG/KG 1,2-Dichloroethane UG/KG Trichloroethene (Trichloroethylene) 12 U 12 U UG/KG Methylcyclohexane 12 U UG/KG 1,2-Dichloropropane 12 U UG/KG Bromodichloromethane UG/KG 12 U cis-1,3-Dichloropropene 12 UJ UG/KG Methyl Isobutyl Ketone 12 U UG/KG Toluene UG/KG 12 U trans-1,3-Dichloropropene 12 U UG/KG 1.1.2-Trichloroethane 12 U UG/KG Tetrachloroethene (Tetrachloroethylene) 12 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

13 UJ UG/KG

Methyl Butyl Ketone

Sample	6796 F	Y 2003 Project: 03-047	4				Produ	iced by: Goddard, Do	enise		
Volatiles	Scan						Requ				•
-						٠	Proje	ot Leader: BSTRIGG	0	)	
_		Creosoting Flowd	ood, MS				Begin	ning: 04/22/2003 16	:50		
Program:	SF		Case No: 31635				Endin	a:			•
Id/Station:	GS06SB /	:	MD No: 1XX6	Inora C	ontractor: S	ENTIN		3.			
		E SOIL (> 12")	D No: 1XX6		ntractor: LIE		DATA	REPORTED ON DE	RY WEIGHT B	ASIS	
RESULTS	UNITS	ANALYTE		<del></del>		RESULTS	UNITS	ANALYTE			
13 U	UG/KG	Dichlorodifluoromethane				13 U	UG/KG	Dibromochlorome	thane		•
13 U	UG/KG	Chloromethane	•	•		13 U	UG/KG	1,2-Dibromoethan			
13 U	UG/KG	Vinyl Chloride				13 U	UG/KG	Chlorobenzene	C (LDD)		•
.13 U	UG/KG	Bromomethane	•			13 U	UG/KG	Ethyl Benzene			
13 U	UG/KG	Chloroethane			•	13 U	UG/KG	Total Xylenes			•
13 U	UG/KG	Trichlorofluoromethane (F	reon 11)			13 U	UG/KG	Styrene			
13 U	UG/KG	1,1-Dichloroethene (1,1-Di				13 U	UG/KG	Bromoform			
13 U	UG/KG	1,1,2-Trichloro-1,2,2-Triflu	oroethane (Freon 113)			13 U	UG/KG	Isopropylbenzene	•		
21 J	UG/KG	Acetone	ordemane (r redir 110)			13 U	UG/KG	1,1,2,2-Tetrachlor			
13 U	* UG/KG	Carbon Disulfide				13 U	UG/KG	1,3-Dichlorobenze			•
13 U	UG/KG	Methyl Acetate				13 U	UG/KG	1,4-Dichlorobenze			•
13 U.	UG/KG	Methylene Chloride				13 U	UG/KG	1,2-Dichlorobenze			
13 U	UG/KG	trans-1,2-Dichloroethene				13 U	UG/KG	1,2-Dibromo-3-Ch		BCP)	
13 U.	UG/KG	Methyl T-Butyl Ether (MTB	E)			13 U	UG/KG	1,2,4-Trichloroben		,,	
13 U	UG/KG	1,1-Dichloroethane	_,			NA	UG/KG	1,2,3-Trichloroben			•
13 U	UG/KG	cis-1,2-Dichloroethene				26	%	% Moisture			
13 UJ	UG/KG	Methyl Ethyl Ketone						,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,			
NA	UG/KG	Bromochloromethane				•			•		
13 U	UG/KG	Chloroform							•		
13 U	UG/KG	1,1,1-Trichloroethane						•		•	
13 Ū	UG/KG	Cyclohexane		•							
13 U	UG/KG	Carbon Tetrachloride	•								
13 U	UĠ/KG	Benzene									
13 U	UG/KG	1,2-Dichloroethane								•	
13 U	UG/KG	Trichloroethene (Trichloroe	ethylene)	•				•			
13 U	UG/KG	Methylcyclohexane	,								
13 U	UG/KG	1,2-Dichloropropane	, ^		•		•				
13 U	UG/KG	Bromodichloromethane			•						
. 13 U	UG/KG	cis-1,3-Dichloropropene			•				•		
13 UJ	UG/KG	Methyl Isobutyl Ketone									
13 U	UG/KG	Toluene	•								
13 U	UG/KG	trans-1,3-Dichloropropene									
13 U	UG/KG	1,1,2-Trichloroethane								•	
13 U	UG/KG	Tetrachloroethene (Tetrach	nloroethylene)					* .			
								· ·			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**VOLATILES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6796 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 16:50

Ending:

**MISCELLANEOUS COMPOUNDS** Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS06SB /

MD No: 1XX6

D No: 1XX6

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

Org Contractor: LIBRTY

**RESULTS UNITS** 

**ANALYTE** 

UG/KG

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

UG/KG

Methyl Butyl Ketone

13 UJ

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample \_ **6806** Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/23/2003 08:45 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS07SB / MD No: 1XY6 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XY6 DATA REPORTED ON DRY WEIGHT BASIS

Media: SUBSURFACE SUIL (> 12 )		DINO: TXYO OIG CONTRACTOR. LIBRET			DATA REPORTED ON DRY WEIGHT BASIS						
RESULTS	UNITS	ANALYTE					RESULT	S UNITS	ANALYTE		
13 U	UG/KG	Dichlorodifluoromethane					13 U	UG/KG	Dibromochloromethane	•	
13 U	UG/KG	Chloromethane					13 U	UG/KG	1,2-Dibromoethane (EDB)		
13 U	UG/KG	Vinyl Chloride					13 U	UG/KG	Chlorobenzene	,	
13 UJ	UG/KG	Bromomethane	•				13 U	UG/KG	Ethyl Benzene		•
13 U	UG/KG	Chloroethane					13 U	UG/KG	Total Xylenes		
13 U	UG/KG	Trichlorofluoromethane	(Freon 11)				13 U	UG/KG	Styrene		
13 U	UG/KG	1,1-Dichloroethene (1,1	Dichloroethylene)				13 U	UG/KG	Bromoform		
13 U	UG/KG	1,1,2-Trichloro-1,2,2-Tril	luoroethane (Freor	n 113)			13 U	UG/KG	Isopropylbenzene		
13 UJ	UG/KG	Acetone					13 U	UG/KG	1,1,2,2-Tetrachloroethane		
13 U	UG/KG	Carbon Disulfide					- 13 U	UG/KG	1,3-Dichlorobenzene		
13 U	UG/KG	Methyl Acetate					13 U	UG/KG	1,4-Dichlorobenzene		
13 Ų	UG/KG	Methylene Chloride					13 U	UG/KG	1,2-Dichlorobenzene		
13 U	UG/KG	trans-1,2-Dichloroethene					13 UJ		1,2-Dibromo-3-Chloropropar	ie (DBCP)	
13 UJ	UG/KG	Methyl T-Butyl Ether (M)	TBE)				13 U	UG/KG	1,2,4-Trichlorobenzene		
13 U	UG/KG	1,1-Dichloroethane					NA		1,2,3-Trichlorobenzene		•
13 U	UG/KG	cis-1,2-Dichloroethene					26	%	% Moisture		
13 UJ	UG/KG	Methyl Ethyl Ketone			•						,
NA	UG/KG	Bromochloromethane		•		•					
	- UG/KG	Chloroform									•
13 U	UG/KG	1,1,1-Trichloroethane							•	•	·
13 U	UG/KG	Cyclohexane								•	
13 Ų	UG/KG	Carbon Tetrachloride	٠,						•		
13 U	UG/KG	Benzene									
13 U	UG/KG	1,2-Dichloroethane									
13 U	UG/KG	Trichloroethene (Trichlor	oethylene)				•				
13 U	ÜG/KG	Methylcyclohexane									
13 U	UG/KG	1,2-Dichloropropane									
13 U	UG/KG	Bromodichloromethane	•						•		
13 U	UG/KG	cis-1,3-Dichloropropene									
13 UJ	UG/KG	Methyl Isobutyl Ketone			•						
13 U	UG/KG	Toluene							•	•	
13 U	UG/KG	trans-1,3-Dichloroproper	ie				-				
13 U	UG/KG	1,1,2-Trichloroethane	-1-1			•					
13 U	UG/KG	Tetrachloroethene (Tetra	icnioroetnylene)						•		•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

	Sample	<b>6804</b> F	Y <b>2003</b> Pr	oject: <b>03-047</b>	<b>'</b> 4					luced by: Goddard, Denise			
	Volatiles		•							uestor:			
			, .	:	•				Proje	ect Leader: BSTRIGGO	•		
			Creosoting	Flow	ood, MS				Begi	nning: 04/23/2003 08:20			
	Program:	SF			Case No: 31635	•			Endi	ng:			
	ld/Station:	GS08SB /			MD No: 1XY4	Inorg Contrac	ctor: SE	ENTIN					
			E SOIL (> 12")	۱ ا	D No: 1XY4	Org Contracto	or: LIB	RTY	DAT	A REPORTED ON DRY WEIG	HT BASIS		
<u>.                                    </u>			2 0012 (> 12	′						A NEI ONTED ON DITT WEIG		· · · · · · · · · · · · · · · · · · ·	_
	RESULTS	UNITS	ANALYTE			•	•	RESULTS	UNITS	ANALYTE	-		
	10 U	UG/KG	Dichlorodiflu	oromethane				10 U	UG/KG	Dibromochloromethane		•	
	10 U	UG/KG	Chlorometh	ane				10 U	UG/KG	1,2-Dibromoethane (EDB)			
	- 10 U	UG/KG	Vinyl Chloric	de :	. •			10 U	UG/KG	Chlorobenzene			
	10 UJ	UG/KG	Bromometh	ane				10 U	UG/KG	Ethyl Benzene			
	10 U	UG/KG	Chloroethan					10 U	UG/KG	Total Xylenes		•	
	10 U	UG/KG		oromethane ((l				10 U	UG/KG	Styrene	•	•	
	10 U	UG/KG	1,1-Dichloro	ethene (1,1-D	ichloroethylene)			10 U	UG/KG	Bromoform			
	ָ 10 U	UG/KG		pro-1,2,2-Triflu	ioroethane (Freon 113)			10 U	UG/KG	Isopropylbenzene			
	. 99 J	UG/KG	Acetone	:				10 U	UG/KG	1,1,2,2-Tetrachloroethane			
		. UG/KG	Carbon Disc					10 U	UG/KG	1,3-Dichlorobenzene			
	10 U	UG/KG	Methyl Acet					10 U	UG/KG	1,4-Dichlorobenzene			
	10 U	UG/KG	Methylene C	Chloride				10 U	UG/KG	1,2-Dichlorobenzene			
	10 U	UG/KG		chloroethene				10 UJ	UG/KG	1,2-Dibromo-3-Chloropropa	ane (DBCP)		
٠.	10 UJ	UG/KG		tyl Ether (MTE	3⊨)			10 U	UG/KG	1,2,4-Trichlorobenzene		,	
	10 U	UG/KG UG/KG	1,1-Dichloro						UG/KG	1,2,3-Trichlorobenzene		-	•
	10 U	UG/KG UG/KG	cis-1,2-Dich			•		17	%	% Moisture		•	
	11 J NA	UG/KG UG/KG	Methyl Ethyl Bromochloro				-						
	10 U	UG/KG	Chloroform	omethane									
	10 U	UG/KG	1,1,1-Trichle	roothana						i			
	10 U	UG/KG	Cyclohexane		•								
	10 U	UG/KG	Carbon Tetra										
	10 U	UG/KG	Benzene	acilionae	* - * - *								
	10 U	UG/KG	1,2-Dichloro	ethane	•								
	10 U	UG/KG		ene (Trichloro	ethylene)			•		•			
	10 U	UG/KG	Methylcyclor										
	10 U	UG/KG	1,2-Dichloro		•								
	10 U	UG/KG	Bromodichlo										
	10 U	UG/KG	cis-1,3-Dichl	oropropene	•			•		•			
	10 UJ	UG/KG	Methyl Isobu			•						•	
	10 U	UG/KG	Toluene		•					•			
	10 U	UG/KG		chloropropene	•								
	10 U	UG/KG	1,1,2-Trichlo			•							
	10 U	UG/KG	Tetrachloroe	thene (Tetrac	hloroethylene)								
	10 UJ	UG/KG	Methyl Butyl	Ketone									
				į.	•								
					•								

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**VOLATILES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6804 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Flowood, MS Facility: Gulf States Creosoting

Program: SF

Project Leader: BSTRIGGO

Beginning: 04/23/2003 08:20

Id/Station: GS08SB /

Case No: 31635

Ending:

Media: SUBSURFACE SOIL (> 12")

**MISCELLANEOUS COMPOUNDS** 

MD No: 1XY4 D No: 1XY4

Inorg Contractor: SENTIN Org Contractor: LIBRTY

RESULTS. UNITS

**ANALYTE** 

8 J - UG/KG

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

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11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 U

11 UJ

11 UJ

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

1,2-Dichloroethane

Methylcyclohexane

1,2-Dichloropropane

Bromodichloromethane

cis-1.3-Dichloropropene

trans-1,3-Dichloropropene

Methyl Isobutyl Ketone

1,1,2-Trichloroethane

Methyl Butyl Ketone

Toluene

Trichloroethene (Trichloroethylene)

Tetrachloroethene (Tetrachloroethylene)

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Sample 6791 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:25 Program: SF Case No: 31635 Endina: Id/Station: GS09SB / MD No: 1XX1 Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")	D No: 1XX1 Org Contractor: LIBRTY			DATA REPORTED ON DRY WEIGHT BASIS				
RESULTS UNITS ANALYTE	:		RESULTS	UNITS	ANALYTE			
11 U UG/KG Dichlorodifluorometh	ane <sup>:</sup>		11 U	UG/KG	Dibromochloromethane	• • •		
11 U UG/KG Chloromethane	;		11 U	UG/KG	1,2-Dibromoethane (EDB)			
11 U UG/KG Vinyl Chloride	:		11 U	UG/KG	Chlorobenzene			
11 U UG/KG Bromomethane			11 U	UG/KG	Ethyl Benzene			
11 U UG/KG Chloroethane		-	11 U	UG/KG	Total Xylenes	•		
11 U UG/KG Trichlorofluorometha	ne (Freon 11)		11 U	UG/KG	Styrene			
11 U UG/KG 1,1-Dichloroethene (	1,1-Dichloroethylene)		11 U	UG/KG	Bromoform	•		
11 U UG/KG 1,1,2-Trichloro-1,2,2	Trifluoroethane (Freon 113)		11 U	UG/KG	Isopropylbenzene			
11 UJ UG/KG Acetone			11 U	UG/KG .	1,1,2,2-Tetrachloroethane			
11 U UG/KG Carbon Disulfide	•	,	11 U	UG/KG	1,3-Dichlorobenzene	•		
11 U UG/KG Methyl Acetate			11 U	UG/KG	1,4-Dichlorobenzene			
11 U UG/KG Methylene Chloride	:		11 U	UG/KG	1,2-Dichlorobenzene			
11 U UG/KG trans-1,2-Dichloroeth			11 U	UG/KG	1,2-Dibromo-3-Chloropropane (D	BCP)		
11 U UG/KG Methyl T-Butyl Ether	(MTBE)		11 U	UG/KG	1,2,4-Trichlorobenzene			
11 U UG/KG 1,1-Dichloroethane		•	NA	UG/KG	1,2,3-Trichlorobenzene			
11 U UG/KG cis-1,2-Dichloroether	ne .		21	%	% Moisture			
11 UJ UG/KG Methyl Ethyl Ketone	1							
NA UG/KG Bromochloromethan	9 .							
. 11 U UG/KG Chloroform								
11 U TUG/KG 1,1,1-Trichloroethane								
11 U UG/KG Cyclohexane								
11 U UG/KG Carbon Tetrachloride						•		
11 U UG/KG Benzene					•			

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample	6786 FY	2003	Project:	03-047	4					Produ	uced by: Goddard, Denise			
Volatiles :			7							Requ	estor:	•		-
			•							Proje	ct Leader: BSTRIGGO			
	oulf States C	rebsoting	)	Flowe	ood, MS					Begir	nning: 04/22/2003 15:15			
Program:				-	Case No: 31635					Endir	ng:			
Id/Station:	GS10SB /				MD No: 1XW6		Inorg Contra	ctor: S	SENTIN					
Media: SU	BSURFACE	SOIL (>	12")	į.	D No: 1XW6		<b>Org Contract</b>	or: Lif	BRTY	DATA	REPORTED ON DRY WEIGH	IT BASIS		
 		<u>`</u>	<u></u>					• •						
RESULTS		ANALY					•		RESULTS	UNITS	ANALYTE			
13.U	UG/KG		difluorom	ethane					13 U	UG/KG	Dibromochloromethane			
13 U	UG/KG	Chlorom							13 U	UG/KG	1,2-Dibromoethane (EDB)			
13 U 13 UJ	UG/KG UG/KG	Vinyl Ch Bromom							13 U 13 U	UG/KG UG/KG	Chlorobenzene Ethyl Benzene			
13 U	UG/KG	Chloroet		İ	*.			•	13 U	UG/KG	Total Xylenes			
13 U	UG/KG		ofluorome	thane /F	reon 11\				13 U	UG/KG	Styrene		•	
13 U	UG/KG				ichloroethylene)				13 U	UG/KG	Bromoform		•	
13 U	UG/KG				oroethane (Freon	113)		•	13 U	UG/KG	Isopropylbenzene			· ·
78 J	UG/KG	Acetone		,		,			13 U	UG/KG	1,1,2,2-Tetrachloroethane			
	. UG/KG	Carbon	Disulfide				•		13 U	UG/KG	1,3-Dichlorobenzene			•
13 U	UG/KG	Methyl A							13 U	UG/KG	1,4-Dichlorobenzene			
13 U	UG/KG		ne Chlorid		•		•		13 U	UG/KG	1,2-Dichlorobenzene			
13 U	UG/KG		2-Dichloro		<del>-</del> \				13 U	UG/KG	1,2-Dibromo-3-Chloropropa	1e (DBCP)		
13 U	UG/KG		-Butyl Eth		E)				13 U	UG/KG	1,2,4-Trichlorobenzene			_
13 U 13 U	UG/KG UG/KG		loroethan Dichloroetl		•					UG/KG %	1,2,3-Trichlorobenzene % Moisture			
13 U	UG/KG	Mothyl F	thyl Ketor	nene	•				17	70	% Woisture			•
NA	UG/KG		lorometh						•					
13 U	UG/KG	Chlorofo		1			•							
13 U	UG/KG		chloroetha	ane	•					,				
13 U	UG/KG	Cyclohe	xane				•	•						
13 U	UG/KG	Carbon	<b>Fetrachlor</b>	ide	•									
2 J	UG/KG	Benzene											•	
13 U	UG/KG		loroethan			•								
13 U	UG/KG		ethene (T		ethylene)	•								
13 U 13 U	UG/KG UG/KG		clohexane loropropa		•		_		•		•			
13 U	UG/KG	Bromodi	chloromet	hane										
13 U	UG/KG		ichloropro			_								
13 U	UG/KG		obutyl Ke								•			
13 U	UG/KG	Toluene	-	-							•			
13,U	UG/KG		-Dichloro									-		
13 U	UG/KG		chloroetha											
13 U	UG/KG				nloroethylene)						•			
13 U	UG/KG	methyl B	utyl Ketor	ne										
	<b>x</b> .													
-				!	•									

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit. | Reporting limit. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**VOLATILES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6786 FY 2003 Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 15:15

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS10SB /

Media: SUBSURFACE SOIL (> 12")

MD No: 1XW6

D No: 1XW6

Inorg Contractor: SENTIN Org Contractor: LIBRTY

**RESULTS UNITS** 25 J UG/KG

ANALYTE

2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise 6789 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:10 Program: SF Case No: 31635 Endina: Id/Station: GS11SB / MD No: 1XW9 Inorg Contractor: SENTIN Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XW9 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS RESULTS UNITS ANALYTE** ANALYTE 11 U UG/KG Dichlorodifluoromethane 11 U UG/KG Dibromochloromethane 11 U UG/KG Chloromethane 11 U UG/KG 1,2-Dibromoethane (EDB) . 11 U UG/KG Vinyl Chloride 11 U UG/KG Chlorobenzene - 11 U UG/KG Bromomethane 11 U Ethyl Benzene UG/KG 11 U UG/KG Chloroethane 11 U UG/KG **Total Xvlenes** 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene 1,1-Dichloroethene (1,1-Dichloroethylene) 11 U UG/KG 11 U UG/KG Bromoform 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U UG/KG 11 U UG/KG Isopropylbenzene 24 J UG/KG Acetone 1,1,2,2-Tetrachloroethane 11.U UG/KG 11 U UG/KG Carbon Disulfide 11 U UG/KG 1,3-Dichlorobenzene 11 U UG/KG Methyl Acetate 11 U UG/KG 1.4-Dichlorobenzene 11 U UG/KG Methylene Chloride 11 U UG/KG 1.2-Dichlorobenzene 11 U UG/KG trans-1,2-Dichloroethene 11 U UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 11 U UG/KG Methyl T-Butyl Ether (MTBE) 11 U UG/KG 1.2.4-Trichlorobenzene 11 U UG/KG 1.1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene 11 U UG/KG cis-1,2-Dichloroethene % Moisture 17 11 UJ UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane UG/KG 11 U Chloroform 11 U . UG/KG 1.1.1-Trichloroethane 11 U. UG/KG Cyclohexane 11 U UG/KG Carbon Tetrachloride 11 U UG/KG Benzene 11 U UG/KG 1,2-Dichloroethane 11 U UG/KG Trichloroethene (Trichlorbethylene) 11 U UG/KG Methylcyclohexane 1.2-Dichloropropane 11 U UG/KG 11 U UG/KG Bromodichloromethane UG/KG 11 U cis-1,3-Dichloropropene 11 UJ UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene UG/KG 11 U trans-1,3-Dichloropropene 11 U UG/KG 1,1,2-Trichloroethane 11 U UG/KG Tetrachloroethene (Tetrachloroethylene) 11 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample 6799 FY 2003 Project: 03-0474 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 07:50 Program: SF Case No: 31635 Endina: Id/Station: GS12SB / Inorg Contractor: SENTIN MD No: 1XY0 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XY0 DATA REPORTED ON DRY WEIGHT BASIS **ANALYTE** RESULTS **RESULTS UNITS** UNITS **ANALYTE** 11 U UG/KG Dichlorodifluoromethane 11 U UG/KG Dibromochloromethane 11 U UG/KG Chloromethane 11 U UG/KG 1.2-Dibromoethane (EDB) Vinyl Chloride 11 U 11 U UG/KG UG/KG Chlorobenzene 11 UJ UG/KG Bromomethane 11 U UG/KG Ethyl Benzene 11 U UG/KG Chloroethane 11 U UG/KG Total Xvienes UG/KG Trichlorofluoromethane (Freon 11) 11 U 11 U UG/KG Styrene UG/KG 1.1-Dichloroethene (1.1-Dichloroethylene) 11 U UG/KG 11 U Bromoform 1.1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U . 11 U UG/KG UG/KG Isopropylbenzene 11 UJ UG/KG Acetone 11 U UG/KG 1,1,2,2-Tetrachloroethane 11. U UG/KG Carbon Disulfide 11 U UG/KG 1.3-Dichlorobenzene 11 U UG/KG Methyl Acetate 11 U UG/KG 1.4-Dichlorobenzene 11 U UG/KG Methylene Chloride 11 U UG/KG 1.2-Dichlorobenzene 11 U UG/KG trans-1,2-Dichloroethene 11 UJ UG/KG 1.2-Dibromo-3-Chloropropane (DBCP) 11 UJ UG/KG Methyl T-Butyl Ether (MTBE) 11 U 1.2.4-Trichlorobenzene UG/KG 11 U UG/KG 1.1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene 11 U UG/KG cis-1.2-Dichloroethene 26 % Moisture 11 UJ UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 11 U-UG/KG Chloroform 11 U UG/KG 1,1,1-Trichloroethane 11 U UG/KG Cyclohexane UG/KG Carbon Tetrachloride 11 U 11 U UG/KG Benzene 11 U UG/KG 1,2-Dichloroethane UG/KG Trichloroethene (Trichloroethylene) 11 U 11 U UG/KG Methylcyclohexane 11 U UG/KG 1.2-Dichloropropane UG/KG Bromodichloromethane 11 U 11 U UG/KG cis-1,3-Dichloropropene 11 UJ UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene 11 U UG/KG trans-1.3-Dichloropropene 11 U UG/KG 1.1.2-Trichloroethane 11 U UG/KG Tetrachloroethene (Tetrachloroethylene) 11 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. | N-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. | N-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. | N-Presumptive evidence analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise 6774 FY 2003 Project: 03-0474 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:25 Program: SF Case No: 31635 Endina: Inorg Contractor: SENTIN Id/Station: GS13SB / MD No: 1XT4 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XT4 DATA REPORTED ON DRY WEIGHT BASIS RESULTS UNITS **RESULTS UNITS ANALYTE** ANALYTE 10 U UG/KG Dichlorodifluoromethane 10 U UG/KG Dibromochloromethane 10 U \_ UG/KG 10 U UG/KG 1,2-Dibromoethane (EDB) Chloromethane 10 U Vinvl Chloride 10 U UG/KG Chlorobenzene UG/KG 10 UJ UG/KG Bromomethane 10 U UG/KG Ethyl Benzene 10 U UG/KG Chloroethane 10 U UG/KG **Total Xylenes** 10 U· UG/KG Trichlorofluoromethane (Freon 11) 10 U UG/KG Styrene 10 U UG/KG 1,1-Dichloroethene (1,1-Dichloroethylene) 10 U UG/KG Bromoform 10 U UG/KG 1,1,2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 10 U UG/KG Isopropylbenzene 1.1.2.2-Tetrachloroethane 21 UJ UG/KG Acetone 10 U UG/KG 10 U Carbon Disulfide 10 U UG/KG UG/KG 1.3-Dichlorobenzene 10 U 10 U UG/KG Methyl Acetate UG/KG 1.4-Dichlorobenzene UG/KG Methylene Chloride 10 U UG/KG 1,2-Dichlorobenzene 10 U 10 U UG/KG trans-1,2-Dichloroethene 10 UJ UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 10 U UG/KG 10 U 1,2,4-Trichlorobenzene Methyl T-Butyl Ether (MTBE) UG/KG 10 U UG/KG 1.1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene cis-1,2-Dichloroethene 10 U UG/KG % Moisture 15 10 UJ UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 10 U UG/KG Chloroform 10 U UG/KG 1.1.1-Trichloroethane 10 U UG/KG Cyclohexane 10 U UG/KG Carbon Tetrachloride Benzene 10 U UG/KG 10 U UG/KG 1,2-Dichloroethane Trichloroethene (Trichloroethylene) 10 U UG/KG 10 U UG/KG Methylcyclohexane 10 U UG/KG 1,2-Dichloropropane 10 U UG/KG Bromodichloromethane 10 U UG/KG cis-1,3-Dichloropropene 10 UJ \* UG/KG Methyl Isobutyl Ketone 10 U UG/KG Toluene 10 U UG/KG trans-1,3-Dichloropropene 1.1.2-Trichloroethane 10 U . UG/KG 10 U UG/KG Tetrachloroethene (Tetrachloroethylene) 10 UJ UG/KG Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyzed in replicate. Reported value is "average" of replicates.

10 U

UG/KG

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise FY 2003 Project: 03-0474 6784 Sample Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 14:50 Program: SF Case No: 31635 Ending: - Id/Station: GS14SB / MD No: 1XW4 Inorg Contractor: SENTIN Media: SUBSURFACE SOIL (> 12") D No: 1XW4 Org Contractor: LIBRTY DATA DEPODTED ON DRY MEIGHT DAGIC R

Media: SUBSURFACE SOIL (> 12")		D No: 1XW4 Org Contractor. LIBH 1 Y			DATA REPORTED ON DRY WEIGHT BASIS					
RESULTS	UNITS	ANALYTE					RESULTS	UNITS	ANALYTE	
10 U	UG/KG	Dichlorodifluoromethane					10 U	UG/KG	Dibromochloromethane	
10 U	UG/KG	Chloromethane					10 U	UG/KG	1,2-Dibromoethane (EDB)	
10 U	UG/KG	Vinyl Chloride					10 U	UG/KG	Chlorobenzene	•
10 UJ	UG/KG	Bromomethane		-			10 U	UG/KG	Ethyl Benzene	
10 U	UG/KG	Chloroethane					10 U	UG/KG	Total Xylenes	
10 U	UG/KG	Trichlorofluoromethane	(Freon 11)				10 U	UG/KG	Styrene	
10 U	UG/KG	1,1-Dichloroethene (1,1-					10 Ū	UG/KG	Bromoform	
10 U	UG/KG		luoroethane (Freon 113)				10 U	UG/KG	Isopropylbenzene	
55 J	UG/KG	Acetone					10 U	UG/KG	1,1,2,2-Tetrachloroethane	
10 U	UG/KG	Carbon Disulfide					10 U	UG/KG	1,3-Dichlorobenzene	
10 U	UG/KG	Methyl Acetate		•			10 U	UG/KG	1,4-Dichlorobenzene	
10 U	UG/KG	Methylene Chloride					10 Ü	UG/KG	1,2-Dichlorobenzene	•
10 U	UG/KG	trans-1,2-Dichloroethene					10 U	.UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)	
10 U	UG/KG	Methyl T-Butyl Ether (M					10 U	UG/KG	1,2,4-Trichlorobenzene	
10 U	UG/KG	1,1-Dichloroethane	•				<sup>*</sup> NA	UG/KG	1,2,3-Trichlorobenzene	
10 U	UG/KG	cis-1,2-Dichloroethene			,		15	%	% Moisture	
10 U	UG/KG	Methyl Ethyl Ketone							·	•
. NA	UG/KG	Bromochloromethane								
10 U	UG/KG	Chloroform		•					•	
10 U	UG/KG	1,1,1-Trichloroethane	•							
10 U `	UG/KG	Cyclohexane								
10 U	UG/KG	Carbon Tetrachloride							· ·	
10 U	UG/KG	Benzene						,		
10 U	UG/KG	1,2-Dichloroethane	•			-	•			
10 U	UG/KG	Trichloroethene (Trichlo	oethylene)						•	
10 U	UG/KG	Methylcyclohexane								
10 U	UG/KG	1,2-Dichloropropane					•		•	
10 U	UG/KG	Bromodichloromethane								
10 U	UG/KG	cis-1,3-Dichloropropene	•						•	
10 U	UG/KG	Methyl Isobutyl Ketone								
10 U	UG/KG	Toluene								•
10 U	UG/KG	trans-1,3-Dichloroproper	ne							
10 U	UG/KG	1,1,2-Trichloroethane								*
10 U	UG/KG	Tetrachloroethene (Tetra	ichloroethylene)							

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

Methyl Butyl Ketone

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**VOLATILES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6784 FY 2003 Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS14SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

MD No: 1XW4

D No: 1XW4

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:50

Produced by: Goddard, Denise

Ending:

Requestor:

RESULTS UNITS 8 J

UG/KG

**ANALYTE** 

UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

6780 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:50 Program: SF Case No: 31635 Ending: ld/Station: GS15SB / MD No: 1XW0 Inorg Contractor: SENTIN Media: SUBSURFACE SOIL (> 12") Org Contractor: LIBRTY D No: 1XW0 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS ANALYTE** 11 U UG/KG 11 U UG/KG Dibromochloromethane Dichlorodifluoromethane 11 U UG/KG Chloromethane 11 U UG/KG 1,2-Dibromoethane (EDB) 11 U Chlorobenzene 11 U \* UG/KG Vinyl Chloride UG/KG 11 UJ UG/KG UG/KG Ethyl Benzene Bromomethane 11 U 11 U UG/KG Chloroethane 11 U UG/KG Total Xylenes 11 U UG/KG Trichlorofluoromethane (Freon 11) 11 U UG/KG Styrene 11 U UG/KG 1.1-Dichloroethene (1.1-Dichloroethylene) 11 U UG/KG Bromoform 11 Ü UG/KG 1.1.2-Trichloro-1,2,2-Trifluoroethane (Freon 113) 11 U Isopropylbenzene UG/KG UG/KG 67 J Acetone 11 U UG/KG 1,1,2,2-Tetrachloroethane 11 U UG/KG Carbon Disulfide 11 U UG/KG 1.3-Dichlorobenzene ...UG/KG 11 U Methyl Acetate 11 U UG/KG 1.4-Dichlorobenzene 11 U UG/KG Methylene Chloride 11 U UG/KG 1,2-Dichlorobenzene 11 U UG/KG trans-1,2-Dichloroethene 11 U UG/KG 1,2-Dibromo-3-Chloropropane (DBCP) 11 U UG/KG Methyl T-Butyl Ether (MTBE) 11·U 1,2,4-Trichlorobenzene UG/KG 11 U UG/KG 1.1-Dichloroethane NA UG/KG 1.2.3-Trichlorobenzene cis-1,2-Dichloroethene 11 U UG/KG % Moisture 21 11 U UG/KG Methyl Ethyl Ketone NA UG/KG Bromochloromethane 11 U UG/KG Chloroform UG/KG 11 U 1,1,1-Trichloroethane 11 U UG/KG Cyclohexane 11 U UG/KG Carbon Tetrachloride UG/KG 11 U Benzene 11 U UG/KG 1,2-Dichloroethane 11 U UG/KG Trichloroethene (Trichloroethylene) 11 U UG/KG Methylcyclohexane 11 U UG/KG 1.2-Dichloropropane 11 U UG/KG Bromodichloromethane .11 U UG/KG cis-1,3-Dichloropropene 11 U UG/KG Methyl Isobutyl Ketone 11 U UG/KG Toluene 11 U UG/KG trans-1.3-Dichloropropene 11 U UG/KG 1.1.2-Trichloroethane 11 U UG/KG Tetrachloroethene (Tetrachloroethylene) UG/KG 11 U Methyl Butyl Ketone

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**VOLATILES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Project: 03-0474

6780 FY 2003

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS15SB /

MD No: 1XW0

Media: SUBSURFACE SOIL (> 12")

D No: 1XW0

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:50

Ending:

**RESULTS UNITS** 20 J

UG/KG

**ANALYTE** 

2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

11 U

11 U

11 U 11 UJ

11 U

11 U

11 U

11 U 11 UJ UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

1,2-Dichloropropane

Bromodichloromethane

cis-1,3-Dichloropropene

trans-1,3-Dichloropropene

Tetrachloroethene (Tetrachloroethylene)

Methyl Isobutyl Ketone

1,1,2-Trichloroethane

Methyl Butyl Ketone

Toluene

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6767 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 11:40 Program: SF Case No: 31635 Ending: Id/Station: GS16SB / MD No: 1XS7 Inorg Contractor: SENTIN Media: SUBSUBEACE SOIL (> 12") D No: 1YS7 Org Contractor: LIBRTY DATA DEDODTED ON DOVINCIOUT DACIS

Media: SUBSURFACE SOIL (> 12")			D No: 1X57 Org Contractor: LIBR11			DATA REPORTED ON DRY WEIGHT BASIS				
RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE			
11 U	UG/KG	Dichlorodifluoromethane			11 U	UG/KG	Dibromochloromethane			
11 Ū	UG/KG	Chloromethane			11 U	UG/KG	1,2-Dibromoethane (EDB)	•		
11 U	UG/KG	Vinyl Chloride		•	11 U	UG/KG	Chlorobenzene	•		
11 ÚJ	UG/KG	Bromomethane			11 U	UG/KG	Ethyl Benzene	•		
11 U	. UG/KG	Chloroethane	•	•	11 U	UG/KG	Total Xylenes			
11 U	UG/KG	Trichlorofluoromethane	(Freon 11)		11 U	UG/KG	Styrene			
11 U	UG/KG	1,1-Dichloroethene (1,1-			. 11 U	UG/KG	Bromoform			
11 U	UG/KG	1,1,2-Trichloro-1,2,2-Triff			11 U	UG/KG	Isopropylbenzene			
11 UJ	UG/KG	Acetone		•	11 U	UG/KG	1,1,2,2-Tetrachloroethane	•		
11 U	UG/KG	Carbon Disulfide			11 U	UG/KG	1,3-Dichlorobenzene			
11 U	UG/KG	Methyl Acetate			11 U	UG/KG	1,4-Dichlorobenzene			
11 U	UG/KG	Methylene Chloride			11 U	UG/KG	1,2-Dichlorobenzene			
11 U	UG/KG	trans-1,2-Dichloroethene			11 UJ	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)	,		
11 U	UG/KG	Methyl T-Butyl Ether (MT	BE)		11 U	UG/KG	1,2,4-Trichlorobenzene			
11 U	UG/KG	1,1-Dichloroethane	1	•	NA	UG/KG	1,2,3-Trichlorobenzene	•		
11 U	UG/KG	cis-1,2-Dichloroethene			22	%	% Moisture			
11 UJ	UG/KG	Methyl Ethyl Ketone	·				·			
NA	UG/KG	Bromochloromethane					•			
11 U	UG/KG	Chloroform		¥.						
11 U	UG/KG	1,1,1-Trichloroethane								
11 U	UG/KG	Cyclohexane								
11 U	UG/KG	Carbon Tetrachloride								
11 U	UG/KG	Benzene								
11 U	UG/KG	1,2-Dichloroethane				-				
11 U	UG/KG	Trichloroethene (Trichlor	pethylene)	•						
11 U	UG/KG	Methylcyclohexane								

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

VAL ATIL		ARADI E	 VOIC
V/ 11 / 11   11			4 - 1 -
VOLATIL	LO O	71811	

13 U

13 U

UG/KG

UG/KG

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6761 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 10:20 Program: SF Case No: 31635 Ending: Id/Station: GS17SB / MD No: 1XS1 Inorg Contractor: SENTIN Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XS1 DATA REPORTED ON DRY WEIGHT BASIS.

ivicula. 30	BSOI II AO	2 301L (> 12 )	D 140. 1X01	Org Contractor.	LIBITIT	DATA	REPORTED ON DRY WEIG	ITTI BASIS	•
RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE		
13 U	UG/KG	Dichlorodifluoromethane			13 U	UG/KG	Dibromochloromethane		•
13 U	UG/KG	Chloromethane	ļ		. 13 U	UG/KG	1,2-Dibromoethane (EDB)		
13 U	UG/KG	Vinyl Chloride			13 U	UG/KG	Chlorobenzene		
ָ13 UJ	* UG/KG	Bromomethane			13 U	UG/KG	Ethyl Benzene		
13 U	UG/KG	Chloroethane			13 U	UG/KG	Total Xylenes		
13 U	UG/KG	Trichlorofluoromethane			13 U	UG/KG	Styrene		•
13 U	UG/KG	1,1-Dichloroethene (1,1-			13 U	UG/KG	Bromoform		
13 U	UG/KG	1,1,2-Trichloro-1,2,2-Trif	luoroethane (Freon 113)		13 U	UG/KG	Isopropylbenzene		
13 U	UG/KG	Acetone			13 U	UG/KG	1,1,2,2-Tetrachloroethane		
13 U	UG/KG	Carbon Disulfide			13 U	UG/KG	1,3-Dichlorobenzene		
13 U	UG/KG	Methyl Acetate			. 13 U	UG/KG	1,4-Dichlorobenzene		
13 U	UG/KG	Methylene Chloride	] 1		13 U	UG/KG	1,2-Dichlorobenzene		
13 U	UG/KG	trans-1,2-Dichloroethene			13 U	UG/KG	1,2-Dibromo-3-Chloropropa	ane (DBCP)	
13 U	UG/KG	Methyl T-Butyl Ether (MT	BE)		13 U	UG/KG	1,2,4-Trichlorobenzene		
13 U	UG/KG	1,1-Dichloroethane			NA.	UG/KG	1,2,3-Trichlorobenzene		
13 U	UG/KG	cis-1,2-Dichloroethene			22	%	% Moisture	• .	
13 U	UG/KG	Methyl Ethyl Ketone							
· NA	UG/KG	Bromochloromethane						•	
13 U	UG/KG UG/KG	Chloroform					•		
13 U 13 U	UG/KG	1,1,1-Trichloroethane Cyclohexane				•		•	•
13 U	UG/KG	Carbon Tetrachloride							
13 U	UG/KG	Benzene.							
13 U	UG/KG	1.2-Dichloroethane				•	·	•	
13 U	UG/KG	Trichloroethene (Trichloro	nethylene)	•	•		•		
13 Ŭ	UG/KG	Methylcyclohexane			•				
13 U	UG/KG	1,2-Dichloropropane			•		·		
13 U	UG/KG	Bromodichloromethane							•
13 U	UG/KG	cis-1,3-Dichloropropene							
13 Ú	UG/KG	Methyl Isobutyl Ketone			•				
	.UG/KG	Toluene							
13 U	UG/KG	trans-1,3-Dichloropropen	e						
- 13 U	UG/KG	1.1.2-Trichloroethane	`			•			
. 5 .7.								•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyzed in replicate. Reported value is "average" of replicates.

Tetrachloroethene (Tetrachloroethylene)

Methyl Butyl Ketone

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample : 6755 FY 2003

**Volatiles Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS18SB /

MD No: 1XR5

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

Org Contractor: LIBRTY D No: 1XR5

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:50

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

			<u> </u>		<del></del>		
RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE
11 U	UG/KG	Dichlorodifluoromethane			11 U	UG/KG	Dibromochloromethane
11 U	UG/KG	Chloromethane			11 U	UG/KG	1,2-Dibromoethane (EDB)
11 U	UG/KG	Vinyl Chloride			11 U	UG/KG	Chlorobenzene
11 UJ	UG/KG	Bromomethane			11 U	UG/KG	Ethyl Benzene
11 U	UG/KG	Chloroethane			11 U	UG/KG	Total Xylenes
11 U	UG/KG	Trichlorofluoromethane	(Freon 11)		- 11 U	UG/KG	Styrene
11 U	UG/KG	1,1-Dichloroethene (1,1-			11 U	UG/KG	Bromoform
11 U	UG/KG		luoroethane (Freon 113)		11 U	UG/KG	Isopropylbenzene
14 UJ	UG/KG	Acetone		•	11 U	UG/KG	1,1,2,2-Tetrachloroethane
11 U	UG/KG	Carbon Disulfide	·		11 U	UG/KG	1,3-Dichlorobenzene
11 U	UG/KG	Methyl Acetate			11 U	UG/KG	1,4-Dichlorobenzene
11 U	UG/KG	Methylene Chloride		•	11 U	UG/KG	1,2-Dichlorobenzene
11 U	UG/KG	trans-1,2-Dichloroethene	ė		11 UJ	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)
11 U	UG/KG	Methyl T-Butyl Ether (M7	BE)		11 U	UG/KG	1,2,4-Trichlorobenzene
11 U	UG/KG	1,1-Dichloroethane	i .		NA	UG/KG	1,2,3-Trichlorobenzene
11 U	UG/KG	cis-1,2-Dichloroethene			25	%	% Moisture
11 UJ	UG/KG	Methyl Ethyl Ketone					
· NA	UG/KG	Bromochloromethane					,
11 U	UG/KG	Chloroform					•
11 U	UG/KG	1,1,1-Trichloroethane					·
11 U	UG/KG	Cyclohexane	1 1	•			
	.UG/KG	Carbon Tetrachloride		•			
11 U	UG/KG	Benzene					
11 U	UG/KG	1,2-Dichloroethane					
· 11 U	UG/KG	Trichloroethene (Trichlor	pethylene)				
11 U	UG/KG	Methylcyclohexane	!				
11 U	UG/KG	.1,2-Dichloropropane	-				
11 U	UG/KG	Bromodichloromethane					
11 U	UG/KG	cis-1,3-Dichloropropene					
11 UJ	UG/KG	Methyl Isobutyl Ketone	!				•
11 U	UG/KG	Toluene	İ				
11 U	UG/KG	trans-1,3-Dichloropropen	e .	•			
11 U	UG/KG	1,1,2-Trichloroethane					•
11 U	UG/KG	Tetrachloroethene (Tetra	chloroethylene)				
11 UJ	UG/KG	Methyl Butyl Ketone					

U-Analyte not detected at or above reporting limit. | J-IdentIfication of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Produced by: Goddard, Denise Sample FY 2003 Project: 03-0474 6757 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 09:35 Program: SF Case No: 31635 Ending: Id/Station: GS19SB / Inorg Contractor: SENTIN MD No: 1XR7 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XR7 DATA REPORTED ON DRY WEIGHT BASIS

				•	J -			טאואט	TEL OTTED ON DITT WEIGHT DAGIO
RESULTS	UNITS	ANALYTE	·				RESULTS	UNITS	ANALYTE
11 U	UG/KG	Dichlorodifluoromethane					11 U	UG/KG	Dibromochloromethane
11 U	UG/KG	Chloromethane					11 U	UG/KG	1,2-Dibromoethane (EDB)
11 U	UG/KG	Vinyl Chloride					11 U	UG/KG	Chlorobenzene
11 UJ	UG/KG	Bromomethane					11 U	UG/KG	Ethyl Benzene
11 U	UG/KG	Chloroethane .					11 U	UG/KG	Total Xylenes
11 U	UG/KG	Trichlorofluoromethane					11 U	UG/KG	Styrene
11 U	UG/KG	1,1-Dichloroethene (1,1-					11 U	UG/KG	Bromoform
11 U	UG/KG	1,1,2-Trichloro-1,2,2-Trif	luoroethane (Freor	า 113)			11 U	UG/KG	Isopropylbenzene
26 UJ	UG/KG	Acetone					11 U	UG/KG	1,1,2,2-Tetrachloroethane
11 U	UG/KG	Carbon Disulfide	' I				11 U	UG/KG	1,3-Dichlorobenzene
11 U	UG/KG	Methyl Acetate					11 U	UG/KG	1,4-Dichlorobenzene
11 U	UG/KG	Methylene Chloride					11 U	UG/KG	1,2-Dichlorobenzene
-11 U	UG/KG	trans-1,2-Dichloroethene	,				11 UJ	UG/KG	1,2-Dibromo-3-Chloropropane (DBCP)
11 U	UG/KG	Methyl T-Butyl Ether (MT	BE)				11 U	UG/ĶG	1,2,4-Trichlorobenzene
11 U	UĠ/KG	1,1-Dichloroethane					· NA	UG/KG	1,2,3-Trichlorobenzene
11 U	UG/KG	cis-1,2-Dichloroethene					21	%	% Moisture
.11 UJ	UG/KG	Methyl Ethyl Ketone							
NA	UG/KG	Bromochloromethane						. *	•
11 U	UG/KG	Chloroform							
11 U	UG/KG	1,1,1-Trichloroethane							
11 U	UG/KG	Cyclohexane							
11 U	UG/KG	Carbon Tetrachloride			•				·
11 U	UG/KG	Benzene						•	
11 U	UG/KG	1,2-Dichloroethane							
11 U	UG/KG	Trichloroethene (Trichloro	petnylene)						
11 U -	UG/KG	Methylcyclohexane	•	•		•			
11 U	UG/KG	1,2-Dichloropropane							•
11 U	UG/KG	Bromodichloromethane							•
11 U	UG/KG	cis-1,3-Dichloropropene	,					•	
11 UJ	UG/KG	Methyl Isobutyl Ketone					•		
11 U	UG/KG	Toluene							·
11 U	UG/KG	trans-1,3-Dichloropropen	е ,						•

Data Reported as Identified by CLP Lab - IDs Not Verified

1.1.2-Trichloroethane

Methyl Butyl Ketone

Tetrachloroethene (Tetrachloroethylene)

11 U

11 U

11 UJ

UG/KG

UG/KG

UG/KG

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6763 FY 2003

**Volatiles Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS20SB /

MD No: 1XS3

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

D No: 1XS3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 10:55

Requestor:

Ending:

RESULTS	UNITS	ANALYTE			•	RESULTS	UNITS	ANALYTE	-
. 11 U	UG/KG	Dichlorodifluoromethane				11 U	UG/KG	Dibromochloromethane	
11 U	UG/KG	Chloromethane			•	- 11 U	UG/KG	1,2-Dibromoethane (EDB)	
. 11 U	UG/KG	Vinyl Chloride	,			11 U	UG/KG	Chlorobenzene	
11 UJ	. UG/KG	Bromomethane		-	•	11 U -	UG/KG	Ethyl Benzene	
11 U	UG/KG	Chloroethane				11 U	UG/KG	Total Xylenes	
11 U	UG/KG	Trichlorofluoromethane	(Freon 11)			11 U	UG/KG	Styrene	
11 U	UG/KG	1,1-Dichloroethene (1,1-	Dichloroethylene)			11 U	UG/KG	Bromoform	
11 U	UG/KG	1,1,2-Trichloro-1,2,2-Triff	uoroethane (Freon 11	·3)		11 U	UG/KG	Isopropylbenzene	
11 U	UG/KG	Acetone		,		11 U	UG/KG	1,1,2,2-Tetrachloroethane	
11 Ū	UG/KG	Carbon Disulfide				11 U	UG/KG	1,3-Dichlorobenzene	
11.U	UG/KG	Methyl Acetate				11 U	UG/KG	1,4-Dichlorobenzene	
11 U	UG/KG	Methylene Chloride				11 U	UG/KG	1,2-Dichlorobenzene	
11 U	UG/KG	trans-1,2-Dichloroethene	•			11 U	UG/KG	1,2-Dibromo-3-Chloropropa	ane (DBCP)
11 U	UG/KG	Methyl T-Butyl Ether (MT	BE)		•	11 U	UG/KG	1,2,4-Trichlorobenzene	, ,
- 11 U	UG/KG	1,1-Dichloroethane				NA	UG/KG	1,2,3-Trichlorobenzene	
11 U	UG/KG	cis-1,2-Dichloroethene				18	%	% Moisture	*
.11 U	UG/KG	Methyl Ethyl Ketone					•		•
NA	UG/KG	Bromochloromethane				-			
11 U	UG/KG	Chloroform	•					• .	
11 U	UG/KG	1,1,1-Trichloroethane					-		
11 U	UG/KG	Cyclohexane	,						
11 U	UG/KG	Carbon Tetrachloride	•				•		
11 U	UG/KG	Benzene	•					·	
11 U	UG/KG	1,2-Dichloroethane			•	•			
11 U	UG/KG	Trichloroethene (Trichlor	pethylene)	•				•	
(11 U	UG/KG	Methylcyclohexane							
11 U	UG/KG	1,2-Dichloropropane							
11 U	UG/KG	Bromodichloromethane							•
11 U	UG/KG	cis-1,3-Dichloropropene	٠.						
11 U	UG/KG	Methyl Isobutyl Ketone							
11 U	UG/KG	Toluene					•		
–	'UG/KG	trans-1,3-Dichloropropen	e					•	
- 11 U	UG/KG	1,1,2-Trichloroethane							
11 U	UG/KG	Tetrachloroethene (Tetra	phioroethylene)						
11 U	UG/KG	Methyl Butyl Ketone							

U-Analyte not detected at or above reporting limit. | J-Ident|fication of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6769 Requestor: **Volatiles Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 12:15 Program: SF Case No: 31635 Ending: Id/Station: GS21SB / Inorg Contractor: SENTIN MD No: 1XS9 Org Contractor: LIBRTY

Media: SUBSURFACE SOIL (> 12")	D No: 1XS9	Org Contractor: LIBRTY		DATA REPORTED ON DRY WEIGHT BASIS			
RESULTS UNITS ANALYTE			RESULTS	UNITS	ANALYTE		
11 U UG/KG Dichlorodifluoromethan	$\mathbf{e}_{\mathbf{i}}^{\mathbf{i}}$	•	11 U	UG/KG	Dibromochloromethane		
11 U UG/KG Chloromethane			11 U	UG/KG	1,2-Dibromoethane (EDB)		
11 U UG/KG Vinyl Chloride			11 U	UG/KG	Chlorobenzene	•	
11 UJ UG/KG Bromomethane	i		11 U	UG/KG	Ethyl Benzene		
11 U UG/KG Chloroethane			11 U	UG/KG	Total Xylenes		
11 U UG/KG Trichlorofluoromethane	(Freon 11)		11 U	UG/KG	Styrene		
11 U UG/KG 1,1-Dichloroethene (1,1	-Dichloroethylene)	i	11 U	UG/KG	Bromoform		
11 U UG/KG 1,1,2-Trichloro-1,2,2-Tr	ifluoroethane (Freon 113)		11 U	UG/KG	Isopropylbenzene		
11 UJ UG/KG Acetone	,		11 U	UG/KG	1,1,2,2-Tetrachloroethane		
11 U UG/KG Carbon Disulfide		•	11 U	UG/KG	1,3-Dichlorobenzene		
11 U UG/KG Methyl Acetate			11 U	UG/KG	1,4-Dichlorobenzene		
11 U UG/KG Methylene Chloride			11 U	UG/KG	1,2-Dichlorobenzene		
11 U UG/KG trans-1,2-Dichloroether		•	11 U	UG/KG	1,2-Dibromo-3-Chloropropane (DBC	;P)	
11 U UG/KG Methyl T-Butyl Ether (M	ITBE)	•	11 U	UG/KG	1,2,4-Trichlorobenzene		
11 U UG/KG 1,1-Dichloroethane				UG/KG	1,2,3-Trichlorobenzene		
11 U UG/KG cis-1,2-Dichloroethene			20	%	% Moisture		
11 UJ UG/KG Methyl Ethyl Ketone				•			
NA UG/KG Bromochloromethane		•		-	•		
11 U UG/KG Chloroform						•	
11 U UG/KG 1,1,1-Trichloroethane	i ·	•	•	•	· ·		
11 U UG/KG Cyclohexane							
11 U UG/KG Carbon Tetrachloride							
11 U UG/KG Benzene	į	•		•	•		
11 U UG/KG 1,2-Dichloroethane							
11 U UG/KG Trichloroethene (Trichlo	roethylene)						
11 U UG/KG Methylcyclohexane	ľ						
11 U UG/KG 1,2-Dichloropropane							
11 U UG/KG Bromodichloromethane	• •	•					
11 U UG/KG cis-1,3-Dichloropropene	<b>)</b>						
11 UJ UG/KG Methyl Isobutyl Ketone	İ						
11 U UG/KG Toluene						•	
11 U UG/KG trans-1,3-Dichloroprope	ne						

11 U

11 U 11 UJ UG/KG

UG/KG

UG/KG

1,1,2-Trichloroethane

Methyl Butyl Ketone

Tetrachloroethene (Tetrachloroethylene)

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

## **UNITED STATES ENVIRONMENTAL PROTECTION AGENCY**

## Region 4

Science and Ecosystem Support Division 980 College Station Road Athens, Georgia 30605-2720

## **MEMORANDUM**

Date: 06/20/2003

Subject: Results of EXTRACTABLES Sample Analysis

03-0474

**Gulf States Creosoting** 

Flowood, MS /

From: Goddard, Denise

To: Striggow, Brian

Thru: QA Office

Attached are the results of analysis of samples collected as part of the subject project. If you have any questions, please contact me.

Produced by: Goddard, Denise Project: 03-0474 Sample 6754 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 09:00 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN ld/Station: GS01SD / MD No: 1XR4 Org Contractor: LIBRTY D No: 1XR4 Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE		RESULTS		ANALYTE	
89 J	UG/KG	Benzaldehyde		440 U	UG/KG	Dibenzofuran	
440 U	UG/KG.	Phenol		440 U	UG/KG	2,4-Dinitrotoluene	
440 U	UG/KG	bis(2-Chloroethyl) Ether		440 U	UG/KG	Diethyl Phthalate	
440 U	UG/KG	2-Chlorophenol		440 U	UG/KG	Fluorene	
440 U	UG/KG	2-Methylphenol		440 U	UG/KG	4-Chlorophenyl Phenyl Ether	
440 U	UG/KG	bis(2-Chloroisopropyl) Ether		1100 U	UG/KG	4-Nitroaniline	,
440 U	UG/KG	Acetophenone		1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol	
440 U	UG/KG	(3-and/or 4-)Methylphenol		440 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine	
440 U	UG/KG	n-Nitrosodi-n-Propylamine		NA	UG/KG	1,2,4,5-Tetrachlorobenzene	
440 U	UG/KG	Hexachloroethane		440 U	UG/KG	4-Bromophenyl Phenyl Ether	
440 U	UG/KG	Nitrobenzene		440 U	UG/KG	Hexachlorobenzene (HCB)	
440 U	UG/KG	Isophorone		440 U	UG/KG	Atrazine	
440 U	UG/KG	2-Nitrophenol		1100 U	UG/KG	Pentachlorophenol	
440 U	UG/KG	2,4-Dimethylphenol		440 U	UG/KG	Phenanthrene	
440 U	UG/KG	bis(2-Chloroethoxy)Methane		440 U	UG/KG	Anthracene	
440 U	UG/KG	2,4-Dichlorophenol		440 U	UG/KG	Carbazole	
440 U	UG/KG	Naphthalene		440 U	UG/KG	Di-n-Butylphthalate	
<sup>1</sup> 440 U	UG/KG	4-Chloroaniline		440 U	UG/KG	Fluoranthene	
440 U	UG/KG	Hexachlorobutadiene		440 U	UG/KG	Pyrene	
440 U	UG/KG	Caprolactam		440 UJ	UG/KG	Benzyl Butyl Phthalate	
440 U	UG/KG	4-Chloro-3-Methylphenol		440 U	UG/KG	3,3'-Dichlorobenzidine	
440 U	UG/KG	2-Methylnaphthalene		440 U	UG/KG	Benzo(a)Anthracene	
440 U	UG/KG	Hexachlorocyclopentadiene (HCCP)		440 U	UG/KG	Chrysene	
440 U	UG/KG	2,4,6-Trichlorophenol	•	440 UJ	UG/KG	bis(2-Ethylhexyl) Phthalate	
1100 Ú	UG/KG	2,4,5-Trichlorophenol		440 U	UG/KG	Di-n-Octylphthalate	
440 U	UG/KG	1,1-Biphenyl		440 U	UG/KG	Benzo(b)Fluoranthene	
440 U	UG/KG	2-Chloronaphthalene		440 U	UG/KG	Benzo(k)Fluoranthene	
1100 U	UG/KG	2-Nitroaniline		440 U	UG/KG	Benzo-a-Pyrene	
440 U	UG/KG	Dimethyl Phthalate		440 U	UG/KG	Indeno (1,2,3-cd) Pyrene	
440 U	UG/KG	2,6-Dinitrotoluene		440 U	UG/KG	Dibenzo(a,h)Anthracene	
440 U	⊍G/KG	Acenaphthylene		92 J	UG/KG	Benzo(ghi)Perylene	
	UG/KG	3-Nitroaniline		55	%	% Moisture	
440 U	UG/KG	Acenaphthene		- <del>-</del>	•		
1100 U	UG/KG	2,4-Dinitrophenol			-		
1100 U	UG/KG	4-Nitrophenol		•			
						•	
•		. !					

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA EXTRACTABLES SAMPLE ANALYSIS** 

Production Date: 06/20/2003 14:10

Project: 03-0474 6754 FY 2003 Sample

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS01SD / Media: SEDIMENT

Case No: 31635 MD No: 1XR4

D No: 1XR4

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:00

Ending:

**RESULTS UNITS** ANALYTE

HEXADECANOIC ACID 420 NJ UG/KG

540 NJ UG/KG 1-HEXADECENE

1500 NJ UG/KG ERGOST-5-N-3-OL, (3.BETA.)-

1100 NJ UG/KG **STIGMASTEROL** 

UG/KG STIGMAST-4-EN-3-ONE 730 NJ 23 UNKNOWN COMPOUNDS 23000 J UG/KG

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyzed in replicate. Reported value is "average" of replicates.

EXTRACTABLES SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

FY 2003 6754

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 09:00

Ending:

SPECIFIED TESTS

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

Program: SF Id/Station: GS01SD /

MD No: 1XR4

Inorg Contractor: SENTIN

D No: 1XR4 Media: SEDIMENT

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.6 J	UG/KG	2-Methylnaphthalene
0.6 J	UG/KG	Naphthalene
3	UG/KG	Acenaphthylene
0.5 J	UG/KG	Acenaphthene
0.5 J	UG/KG	Fluorene
2 J	UG/KG	Phenanthrene
3	UG/KG	Anthracene
2	UG/KG	Fluoranthene
2	UG/KG	Pyrene
1 J	UG/KG	Benzo(a)Anthracene
1 J	UG/KG	Chrysene
NA	·UG/KG	Benzo(b)Fluoranthene
NA	UG/KG	Benzo(k)Fluoranthene
2	.UG/KG	Benzo-a-Pyrene
NA	UG/KG	Indeno (1,2,3-cd) Pyrene
1 U	UG/KG	Dibenzo(a,h)Anthracene
NA	UG/KG	Benzo(ghi)Perylene
24 U	UG/KG	Pentachlorophenol

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences... | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. Media: SUBSURFACE SOIL (> 12")

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise Project: 03-0474 6755 FY 2003 Sample Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 08:50 Program: SF Case No: 31635 Ending: Id/Station: GS18SB / MD No: 1XR5 Inorg Contractor: SENTIN

Org Contractor: LIBRTY

		•		•		·		·	
RESULTS	UNITS	ANALYTE				RESULTS	UNITS	ANALYTE	
440 U	UG/KG	Benzaldehyde		• •	•	440 U	UG/KG	Dibenzofuran	
440 U	UG/KG	Phenol				440 U	UG/KG	2,4-Dinitrotoluene	
440 U	UG/KG	bis(2-Chloroethyl) Ether	•		1	440 U	UG/KG	Diethyl Phthalate	
440 Ú	UG/KG	2-Chlorophenol				440 U	UG/KG	Fluorene	
440 U	· UG/KG	2-Methylphenol				440 U	UG/KG	4-Chlorophenyl Phenyl Ether	
440 U	UG/KG	bis(2-Chloroisopropyl) E	ther			1100 U	UG/KG	4-Nitroaniline	
440 U	UG/KG	Acetophenone				1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol	
440 U	UG/KG	(3-and/or 4-)Methylphen	ol ·			440 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine	
440 U	UG/KG	n-Nitrosodi-n-Propylamir				NA	UG/KG	1,2,4,5-Tetrachlorobenzene	•
440 U	UG/KG	Hexachloroethane		•	•	440 U	UG/KG	4-Bromophenyl Phenyl Ether	
440 U	UG/KG	Nitrobenzene		·		440 U	UG/KG	Hexachlorobenzene (HCB)	
440 U	UG/KG	Isophorone			•	440 U	UG/KG	Atrazine	
440 U	UG/KG	2-Nitrophenol				1100 U	UG/KG	Pentachlorophenol	
440 U	UG/KG	2,4-Dimethylphenol				440 U	UG/KG	Phenanthrene	
440 U	UG/KG	bis(2-Chloroethoxy)Meth	ane		•	440 U	UG/KG	Anthracene	
440 U	UG/KG	2,4-Dichlorophenol				440 U	UG/KG	Carbazole	•
440 U	UG/KG	Naphthalene	•			440 U	UG/KG	Di-n-Butylphthalate	
440 U	UG/KG	4-Chloroaniline				440 U	UG/KG	Fluoranthene	
440 U	UG/KG	Hexachlorobutadiene				440 U	UG/KG	Pyrene	
440 U	UG/KG	Caprolactam				440 UJ	UG/KG	Benzyl Butyl Phthalate	
440 U	UG/KG	4-Chloro-3-Methylphenol			·	440 U	UG/KG	3,3'-Dichlorobenzidine	•
440 U	UG/KG	2-Methylnaphthalene				440 U	·UG/KG	Benzo(a)Anthracene	
440 U	UG/KG	Hexachlorocyclopentadie	ne (HCCP)		•	440 U	UG/KG	Chrysene	
440 U	UG/KG	2,4,6-Trichlorophenol		•	•	440 UJ	UG/KG	bis(2-Ethylhexyl) Phthalate	
1100 U	UG/KG	2,4,5-Trichlorophenol			•	440 U	UG/KG	Di-n-Octylphthalate	
440 U	UG/KG	1,1-Biphenyl	•			440 U	UG/KG	Benzo(b)Fluoranthene	
440 U	UG/KG	2-Chloronaphthalene				440 U	UG/KG	Benzo(k)Fluoranthene	
1100 U	ŲG/KG	2-Nitroaniline		•		440 U	UG/KG	Benzo-a-Pyrene	
440 U	UG/KG	Dimethyl Phthalate	•			440 U	UG/KG	Indeno (1,2,3-cd) Pyrene	
440 U	UG/KG	2,6-Dinitrotoluene				440 U	UG/KG	Dibenzo(a,h)Anthracene	
440 U	UG/KG	Acenaphthylene				440 U	UG/KG	Benzo(ghi)Perylene	
1100 U	UG/KG	3-Nitroaniline				25	%	% Moisture	
440 U	UG/KG	Acenaphthene			•				
1100 U	UG/KG	2,4-Dinitrophenol						·	
1100 U	UG/KG	4-Nitrophenol							
		·		•		•			
		1		_				•	

D No: 1XR5

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

EXTRACTABLES SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Program: SF

230 J

450 J

Id/Station: GS18SB /

6755 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

MD No: 1XR5 D No: 1XR5

Media: SUBSURFACE SOIL (> 12")

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:50

Endina:

**RESULTS UNITS ANALYTE** 

120 NJ UG/KG UG/KG UG/KG

1.2-BENZENEDICARBOXYLIC ACID

2 UNKNOWN PHTHALATES 2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6756 FY 2003 Project: 03-0474

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS18SS /

Media: SURFACE SOIL (0" - 12")

Case No: 31635 MD No: 1XR6

Inorg Contractor: SENTIN Org Contractor: LIBRTY D No: 1XR6

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
430 U	UG/KG	Benzaldehyde	430 U	UG/KG	Dibenzofuran
430 U	UG/KG	Phenol	430 U	UG/KG	2,4-Dinitrotoluene
430 U	UG/KG	bis(2-Chloroethyl) Ether	430 U	UG/KG	Diethyl Phthalate
430 U	UG/KG	2-Chlorophenol	430 U	UG/KG	Fluorene
430 U -	UG/KG	2-Methylphenol	430 U	UG/KG	. 4-Chlorophenyl Phenyl Ether
430 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
430 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
430 U	UG/KG	(3-and/or 4-)Methylphenol	430 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
430 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
430 U	UG/KG	Hexachloroethane	430 U	UG/KG	4-Bromophenyl Phenyl Ether
430 U	UG/KG	Nitrobenzene	430 U	UG/KG	Hexachlorobenzene (HCB)
430 U	UG/KG	Isophorone	430 U	UG/KG	Atrazine
430 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
430 U	UG/KG	2,4-Dimethylphenol	430 U	UG/KG	Phenanthrene
430 U	*UG/KG	bis(2-Chloroethoxy)Methane	430 U	UG/KG	Anthracene
430 U	UG/KG	2,4-Dichlorophenol	430 U	UG/KG	Carbazole
430 U	UG/KG	Naphthalene	430 U	UG/KG	Di-n-Butylphthalate
430 U	UG/KG	4-Chloroaniline	430 U	UG/KG	Fluoranthene
430 U	UG/KG	Hexachlorobutadiene	430 U	UG/KG	Pyrene
430 U	UG/KG	Caprolactam	430 UJ	UG/KG	Benzyl Butyl Phthalate
430 U	UG/KG	4-Chloro-3-Methylphenol	430 U	UG/KG	3,3'-Dichlorobenzidine
430 U	UG/KG	2-Methylnaphthalene	430 U	UG/KG	Benzo(a)Anthracene
430 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	430 U	UG/KG	Chrysene
430 U	UG/KG	2,4,6-Trichlorophenol	430 UJ	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	430 U	UG/KG	Di-n-Octylphthalate
430 U	UG/KG.	1,1-Biphenyl	430 U	UG/KG	Benzo(b)Fluoranthene
430 U	UG/KG	2-Chloronaphthalene	430 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	430 U	UG/KG	Benzo-a-Pyrene
430 U	UG/KG	Dimethyl Phthalate	430 U	UG/KG	Indeno (1,2,3-cd) Pyrene
430 U	UG/KG	2,6-Dinitrotoluene	430 U	UG/KG	Dibenzo(a,h)Anthracene
430 U	UG/KG	Acenaphthylene	430 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	23	%	% Moisture
430 U	UG/KG	Acenaphthene	•		•
1100 U	UG/KG	2,4-Dinitrophenol			•
1100 U	UG/KG	4-Nitrophenol		•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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**EXTRACTABLES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6756 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS18SS /

Media: SURFACE SOIL (0" - 12")

MD No: 1XR6

D No: 1XR6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:40

Ending:

**ANALYTE RESULTS UNITS** 

110 NJ UG/KG 3100 J UG/KG HEXADECANOIC ACID

17 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6757 FY 2003

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS19SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635 MD No: 1XR7

D No: 1XR7

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:35

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
420 U	UG/KG	Benzaldehyde	420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	420 U	UG/KG	Diethyl Phthalate
420 U	UG/KG	2-Chlorophenol	420 U	UG/KG	Fluorene .
420 U	UG/KG	2-Methylphenol	420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
420 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
420 U	UG/KG	(3-and/or 4-)Methylphenol	· 420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane	420 U	UG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	420 U	UG/KG	Hexachlorobenzene (HCB)
420 U	UG/KG	Isophorone	420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
420 U	UG/KG	2,4-Dimethylphenol	420 U	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Methane	420 U	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol	420 U	UG/KG	Carbazole
420 U	UG/KG	Naphthalene	420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline	420 U	UG/KG	Fluoranthene
420 U	UG/KG	Hexachlorobutadiene	420 U	UG/KG	Pyrene
420 U .	UG/KG	Caprolactam	420 UJ	UG/KG	Benzyl Butyl Phthalate
420 U	UG/KG	4-Chloro-3-Methylphenol	420 U	UG/KG	3,3'-Dichlorobenzidine
420 U	UG/KG	2-Methylnaphthalene	420 U	UG/KG	Benzo(a)Anthracene
420 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	420 U	UG/KG	Chrysene
420 U	_UG/KG	2,4,6-Trichlorophenol	420 UJ	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	420 U	UG/KG	Di-n-Octylphthalate
420 U	UG/KG	1,1-Biphenyl	420 U	UG/KG	Benzo(b)Fluoranthene
420 U	UG/KG	2-Chloronaphthalene	420 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	420 U	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate	420 U	UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2,6-Dinitrotoluene	420 U	UG/KG	Dibenzo(a,h)Anthracene
420 U	UG/KG	Acenaphthylene	420 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	21	%	% Moisture
420 U	UG/KG	Acenaphthene			
1100 U	UG/KG	2,4-Dinitrophenol			·
1100 U	UG/KG	4-Nitrophenol			

U-Analyte not detected at or above reporting limit. | J-Idenţification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EXTRACTABLES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6757 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 09:35

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

95 J

Case No: 31635

ld/Station: GS19SB /

MD No: 1XR7

Media: SUBSURFACE SOIL (> 12")

D No: 1XR7

Inorg Contractor: SENTIN Org Contractor: LIBRTY

UG/KG

**RESULTS UNITS ANALYTE** 

UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

400 U

400 U

1000 U

400 U

400 U

400 U

400 U

1000 U

1000 U

1000 U

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

1.1-Biphenyl

2-Nitroaniline

2-Chloronaphthalene

Dimethyl Phthalate

2.6-Dinitrotoluene

Acenaphthylene

3-Nitroaniline

4-Nitrophenol

Acenaphthene

2.4-Dinitrophenol

Produced by: Goddard, Denise Sample 6758 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:30 Program: SF Case No: 31635 Ending: Id/Station: GS19SS / Inorg Contractor: SENTIN MD No: 1XR8 D No: 1XR8 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** RESULTS UNITS ANALYTE 400 U UG/KG 400 U UG/KG Dibenzofuran Benzaldehyde 400.U UG/KG 2.4-Dinitrotoluene 400 U UG/KG Phenol 400 U UG/KG bis(2-Chloroethyl) Ether 400 U UG/KG Diethyl Phthalate 400 U 400 U UG/KG 2-Chlorophenol UG/KG Fluorene 400 U 400 U UG/KG 2-Methylphenol UG/KG 4-Chlorophenyl Phenyl Ether 400 U bis(2-Chloroisopropyl) Ether 1000 U UG/KG 4-Nitroaniline UG/KG 400 U 1000 U UG/KG 2-Methyl-4,6-Dinitrophenol UG/KG Acetophenone 400 U 400 U UG/KG (3-and/or 4-)Methylphenol UG/KG n-Nitrosodiphenylamine/Diphenylamine 400 U UG/KG 1.2.4.5-Tetrachlorobenzene UG/KG n-Nitrosodi-n-Propylamine NA Hexachloroethane 400 U UG/KG 4-Bromophenyl Phenyl Ether 400 U UG/KG 400 U UG/KG Nitrobenzene 400 U UG/KG Hexachlorobenzene (HCB) 400 U UG/KG Isophorone 400 U UG/KG Atrazine 400 U 2-Nitrophenol 1000 U UG/KG UG/KG Pentachlorophenol 400 U UG/KG 2.4-Dimethylphenol 400 U UG/KG Phenanthrene 400 U UG/KG bis(2-Chloroethoxy)Methane 400 U UG/KG Anthracene 400 U UG/KG 2.4-Dichlorophenol 400 U UG/KG Carbazole 400 U UG/KG Naphthalene 400 U UG/KG Di-n-Butylphthalate 400 U ÚG/KG 4-Chloroaniline 400 U UG/KG Fluoranthene 400 U UG/KG Hexachlorobutadiene 400 U UG/KG Pvrene 400 U UG/KG Caprolactam 400 UJ UG/KG Benzyl Butyl Phthalate 400 U UG/KG 4-Chloro-3-Methylphenol 400 U UG/KG 3,3'-Dichlorobenzidine 400 U UG/KG 2-Methylnaphthalene 400 U UG/KG Benzo(a)Anthracene 400 Ú UG/KG Hexachlorocyclopentadiene (HCCP) 400 U UG/KG Chrysene 400 U UG/KG 2,4,6-Trichlorophenol 400 U UG/KG bis(2-Ethylhexyl) Phthalate 1000 U UG/KG 2,4,5-Trichlorophenol 400 U UG/KG Di-n-Octylphthalate

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

400 U

400 U

400 U

400 U

400 U

400 U

17

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

%

Benzo(b)Fluoranthene

Benzo(k)Fluoranthene

Indeno (1,2,3-cd) Pyrene

Dibenzo(a,h)Anthracene

Benzo(ahi)Pervlene

Benzo-a-Pyrene

% Moisture

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EXTRACTABLES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6758 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS19SS /

MD No: 1XR8

D No: 1XR8

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12") **ANALYTE**  Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:30

Ending:

**RESULTS UNITS** 160 J

UG/KG UG/KG HEXADECANOIC ACID

**TESTOSTERONE** 

390 NJ 12000 J UG/KG

25 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value,

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

FY 2003 Sample

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:50

Ending:

Facility: Gulf States Creosoting

Flowood, MS

Program: SF Id/Station: GS02SD /

**Extractables Scan** 

Case No: 31635

Media: SEDIMENT D No: 1XR9

Inorg Contractor: SENTIN MD No: 1XR9 Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

				· · · · · · · · · · · · · · · · · · ·		
RESULTS	UNITS	ANALYTE		RESULTS	UNITS	ANALYTE
410 U	UG/KG	Benzaldehyde		410 U	UG/KG	Dibenzofuran
410 U	UG/KG	Phenol		410 U	UG/KG	2,4-Dinitrotoluene
410 U	UG/KG	bis(2-Chloroethyl) Ether		410 U	UG/KG	Diethyl Phthalate
410 U.	UG/KG	2-Chlorophenol	•	410 U	UG/KG	Fluorene
410 U	UG/KG	2-Methylphenol		410 U	UG/KG	4-Chlorophenyl Phenyl Ether
410 U	UG/KG	bis(2-Chloroisopropyl) Ether		1000 U	UG/KG	4-Nitroaniline
410 U	UG/KG	Acetophenone		1000 U	UG/KG	2-Methyl-4,6-Dinitrophenol
410 U	UG/KG	(3-and/or 4-)Methylphenol		410 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
410 U	UG/KG	n-Nitrosodi-n-Propylamine		NA	UG/KG	1,2,4,5-Tetrachlorobenzene
410 U	UG/KG	Hexachloroethane	•	410 U	UG/KG	4-Bromophenyl Phenyl Ether
410 U	UG/KG	Nitrobenzene		410 U	UG/KG	Hexachlorobenzene (HCB)
410 U	UG/KG	Isophorone	:	410 U	UG/KG	Atrazine
410 U	UG/KG	2-Nitrophenol		1000 U	UG/KG	Pentachlorophenol
410 U	UG/KG	2,4-Dimethylphenol	•	410 U	UG/KG	Phenanthrene
410 U	UG/KG	bis(2-Chloroethoxy)Methane		410 U	UG/KG	Anthracene
410 U	UG/KG	2,4-Dichlorophenol		410 U	UG/KG	Carbazole
410 U	UG/KG	Naphthalene		410 U	UG/KG	Di-n-Butylphthalate
410 U	UG/KG	4-Chloroaniline		410 U	UG/KG	Fluoranthene
410 U	UG/KG	Hexachlorobutadiene		410 U	UG/KG	Pyrene
410 U	UG/KG	Caprolactam		410 UJ	UG/KG	Benzyl Butyl Phthalate
410 U	UG/KG	4-Chloro-3-Methylphenol		410 UJ	UG/KG	3,3'-Dichlorobenzidine
410 U	UG/KG	2-Methylnaphthalene		410 U	UG/KG	Benzo(a)Anthracene
410 U	UG/KG	Hexachlorocyclopentadiene (HCCP)		. 410 U	UG/KG	Chrysene
410 U	UG/KG	2,4,6-Trichlorophenol		410 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1000 U	UG/KG	2,4,5-Trichlorophenol		410 U	UG/KG	Di-n-Octylphthalate
410 U	UG/KG	1,1-Biphenyl		410 U	UG/KG	Benzo(b)Fluoranthene
410 U	UG/KG	2-Chloronaphthalene		410 U	UG/KG	Benzo(k)Fluoranthene
1000 U	UG/KG	2-Nitroaniline		410 U	UG/KG	Benzo-a-Pyrene .
410 U	UG/KG	Dimethyl Phthalate		410 U	UG/KG	Indeno (1,2,3-cd) Pyrene
410 U	UG/KG	2,6-Dinitrotoluene		410 U	UG/KG	Dibenzo(a,h)Anthracene
410 U	UG/KG	Acenaphthylene		410 U	UG/KG	Benzo(ghi)Perylene
1000 U -	UG/KG	3-Nitroaniline		52	%	% Moisture
410 U	UG/KG	Acenaphthene			•	·
1000 U	UG/KG	2,4-Dinitrophenol	•	•		
1000 U	UG/KG	4-Nitrophenol				
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U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 09:50

Requestor:

Ending:

Sample

6759 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS02SD /

Case No: 31635 MD No: 1XR9

D No: 1XR9

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Media: SEDIMENT **RESULTS UNITS** 

**ANALYTE** 

UG/KG GAMMA.-SITOSTEROL 1400 NJ

28 UNIDENTIFIED COMPOUNDS 92000 J UG/KG

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

## **EXTRACTABLES SAMPLE ANALYSIS**

22 U

UG/KG

Pentachlorophenol

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6759 Requestor: SPECIFIED TESTS Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:50 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS02SD / MD No: 1XR9 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XR9 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** 0.9 J UG/KG 2-Methylnaphthalene Naphthalene 0.6 J UG/KG UG/KG Acenaphthylene UG/KG Acenaphthene 1 J Fluorene 2 UG/KG 12 UG/KG Phenanthrene 7 UG/KG Anthracene 21 UG/KG Fluoranthene 15 UG/KG Pyrene Benzo(a)Anthracene UG/KG 10 UG/KG Chrysene NA UG/KG Benzo(b)Fluoranthene UG/KG Benzo(k)Fluoranthene Benzo-a-Pyrene UG/KG NA UG/KG Indeno (1,2,3-cd) Pyrene Dibenzo(a.h)Anthracene UG/KG Benzo(ghi)Perylene NA UG/KG

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Media: SURFACE SOIL (0" - 12")

Project: 03-0474 Sample 6760 FY 2003

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 10:05

Ending:

Program: SF

Facility: Gulf States Creosoting

Flowood, MS

Id/Station: GS17SS /

**Extractables Scan** 

Case No: 31635

MD No: 1XS0

D No: 1XS0

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	r :			RESULTS		ANALYTE
420 U	UG/KG	Benzaldehyde	<u>.</u>			420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol			•	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	i .			420 U	UG/KG	Diethyl Phthalate
420 U	UG/KG	2-Chlorophenol				420 U	UG/KG	Fluorene
420 U -	UG/KG	2-Methylphenol				420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U	UG/KG	bis(2-Chloroisopropyl) E	ther	•	•	1100 U	UG/KG	4-Nitroaniline
420 U	· UG/KG	Acetophenone				1100 U	ÚG/KG	2-Methyl-4,6-Dinitrophenol
. 420 U	UG/KG	(3-and/or 4-)Methylphen	ol			420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamii			•	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane				420 U	UG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	!			420 U	UG/KG	Hexachlorobenzene (HCB)
420 U	UG/KG	Isophorone				420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol				1100 U	UG/KG	Pentachlorophenol
420 U	UG/KG	2,4-Dimethylphenol	·	•		250 J	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Meth	ane			380 J	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol				130 J	UG/KG	Carbazole
420 U	UG/KG	Naphthalene				420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline			•	2500	UG/KG	Fluoranthene
420 U	UG/KG	Hexachlorobutadiene	: <b>i</b>			1700	UG/KG	Pyrene
420 U	UG/KG	Caprolactam	<u>.</u>	•		420 UJ	UG/KG	Benzyl Butyl Phthalate
420 U	UG/KG	4-Chloro-3-Methylpheno	· 	•		420 UJ	UG/KG	3,3'-Dichlorobenzidine
420 U	· UG/KG	2-Methylnaphthalene				1000	UG/KG	Benzo(a)Anthracene
420 U	UG/KG	Hexachlorocyclopentadie	ne (HCCP)			1600	UG/KG	Chrysene
420 U	UG/KG	2,4,6-Trichlorophenol	( , , , , , , , , , , , , , , , , , , ,	• .		420 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	!	•		420 U	UG/KG	Di-n-Octylphthalate
420 U	UG/KG	1,1-Biphenyl		* .		2600 J	UG/KG	Benzo(b)Fluoranthene
420 U	UG/KG	2-Chloronaphthalene	i			2700 J	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	í			1000	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate				890	UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2,6-Dinitrotoluene	!		•	330 J	UG/KG	Dibenzo(a,h)Anthracene
270 J	UG/KG	Acenaphthylene				290 J	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	•	•		21	%	% Moisture
420 U	UG/KG	Acenaphthene	1	•		٤١	,,	// Moisture
1100 U	UG/KG	2,4-Dinitrophenol						
1100 U	UG/KG	4-Nitrophenol	i ·					
1100 0	Carita	4 This option of	I	· ·				

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable

Sample 6760 FY 2003 Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS17SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

MD No: 1XS0

Inorg Contractor: SENTIN

Beginning: 04/22/2003 10:05 Ending:

Requestor:

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Org Contractor: LIBRTY D No: 1XS0

RESULTS	UNITS	ANALYTE
210 NJ	UG/KG	NAPHTHALENE, 2-PHENYL-
510 NJ	UG/KG	11H-BENZO [B] FLUORENE
390 NJ	UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
310 NJ	UG/KG	BENZO [B] NAPHTHO [2,1-D] THIOPHENE
310 NJ	UG/KG	CHRYSENE, 5-METHYL-
800 NJ	UG/KG	5,12-NAPHTHACENEDIONE
880 NJ	UG/KG	BENZO [E] PYRENE
650 NJ	UG/KG	CHOLESTEROL
290 NJ	UG/KG	1,2:3,4-DIBENZPYRENE
970 J	UG/KG	3 UNKNOWN PAHS
5900 J	UG/KG	16 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

6761 FY 2003 Project: 03-0474 Sample

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS17SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

MD No: 1XS1

D No: 1XS1

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:20

Ending:.

DATA REPORTED ON DRY WEIGHT BASIS

		· · · · · · · · · · · · · · · · · · ·			· · · · · · · · · · · · · · · · · · ·
RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
420 U	UG/KG	Benzaldehyde	420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	420 U	UG/KG	Diethyl Phthalate
420 U	ÚĠ/KĠ	2-Chlorophenol	420 U	UG/KG	Fluorene
420 U	UG/KG	2-Methylphenol	420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
420 U	UG/KG	Acetophenone	1100 U	UG/KG .	2-Methyl-4,6-Dinitrophenol
420 U	UG/KG	(3-and/or 4-)Methylphenol	420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane	420 U	ÚG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	. 420 U	UG/KG	Hexachlorobenzene (HCB)
420 U	UG/KG	Isophorone	420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
420 U	UG/KG	2,4-Dimethylphenol	420 U	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Methane	420 U	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol	420 U	UG/KG	Carbazole
420 U	UG/KG	Naphthalene	420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline	420 U	UG/KG	Fluoranthene
420 Ú	UG/KG	Hexachlorobutadiene	420 U	UG/KG	Pyrene
420 U	UG/KG	Caprolactam	420 UJ	UG/KG	Benzyl Butyl Phthalate
420 U	·UG/KG	4-Chloro-3-Methylphenol	420 UJ	UG/KG	3,3'-Dichlorobenzidine
420 U	UG/KG	2-Methylnaphthalene	420 U	UG/KG	Benzo(a)Anthracene
420 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	420 U	UG/KG	Chrysene
420 U	UG/KG	2,4,6-Trichlorophenol	420 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	420 U	UG/KG	Di-n-Octylphthalate
420 U	UG/KG	1,1-Biphenyl	.41 J	UG/KG	Benzo(b)Fluoranthene
420 U	UG/KG	2-Chloronaphthalene	43 J	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	420 U	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate	420 U	·UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2,6-Dinitrotoluene	420 U	UG/KG	Dibenzo(a,h)Anthracene
420 U	UG/KG	Acenaphthylene	420 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	22	%	% Moisture
420 U	UG/KG	Acenaphthene			
1100 U	UG/KG	2,4-Dinitrophenol	•		
1100 U	UG/KG	4-Nitrophenol			
		•			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample 6762 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 10:45 Program: SF Case No: 31635 Ending: ld/Station: GS20SS / Inorg Contractor: SENTIN MD No: 1XS2 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XS2 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
420 U	UG/KG	Benzaldehyde	420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	420 U	UG/KG	Diethyl Phthalate
420 U	UG/KG	2-Chlorophenol	420 U	UG/KG	Fluorene
420 U	UG/KG	2-Methylphenol	420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U *	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
420 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
420 U	UG/KG	(3-and/or 4-)Methylphenol	420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamine	. NA	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane	420 U	UG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	420 U	UG/KG	Hexachlorobenzene (HCB)
420 U	UG/KG	Isophorone	420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
420 U	UG/KG	2,4-Dimethylphenol	420 U	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Methane	420 U	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol	420 U	UG/KG	Carbazole
420 U	UG/KG	Naphthalene	420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline '	73 J	UG/KG	Fluoranthene
420 U	UG/KG	Hexachlorobutadiene	100 J	UG/KG	Pyrene
420 U	UG/KG	Caprolactam	420 UJ	UG/KG	Benzyl Butyl Phthalate
420 U	UG/KG	4-Chloro-3-Methylphenol	420 UJ	UG/KG	3,3'-Dichlorobenzidine
420 Ų	UG/KG	2-Methylnaphthalene	58 J	UG/KG	Benzo(a)Anthracene
420 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	98 J	UG/KG	Chrysene
420 U	UG/KG	2,4,6-Trichlorophenol	420 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	420 U	UG/KG	Di-n-Octylphthalate .
420 U	UG/KG	1,1-Biphenyl	160 J	UG/KG	Benzo(b)Fluoranthene
420 U	UG/KG	2-Chloronaphthalene	170 J	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	46 J	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate	63 J	UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2,6-Dinitrotoluene	420 U	UG/KG	Dibenzo(a,h)Anthracene
420 U	UG/KG	Acenaphthylene	420 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	21	%	% Moisture
	⊍G/KG	Acenaphthene			•
1100 U	UG/KG	2,4-Dinitrophenol	•		
1100 U	UG/KG	4-Nitrophenol			
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U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

EPA - REGION IV SESD, ATHENS, GA

**EXTRACTABLES SAMPLE ANALYSIS** 

Production Date: 06/20/2003 14:10

Sample 6762 FY 2003 Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

8800 J

. . . . .

Id/Station: GS20SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

MD No: 1XS2

D No: 1XS2

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:45

Ending:

RESULTS UNITS ANALYTE

UG/KG

130 NJ UG/KG 190 NJ UG/KG HEXADECANOIC ACID

1-EICOSANOL

17 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample 6763 FY 2003 Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 10:55

Ending:

**Extractables Scan** 

Facility: Gulf States Creosoting

Media: SUBSURFACE SOIL (> 12")

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS20SB /

MD No: 1XS3

D No: 1XS3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE
400 U	UG/KG	Benzaldehyde			400 U	UG/KG	Dibenzofuran
400 U	UG/KG	Phenol			400 U	UG/KG	2,4-Dinitrotoluene
400 U	UG/KG	bis(2-Chloroethyl) Ether	•		400 U	UG/KG	Diethyl Phthalate
400 U	UG/KG	2-Chlorophenol			400 U	UG/KG	Fluorene
400 U	UG/KG	2-Methylphenol			400 U	UG/KG	4-Chlorophenyl Phenyl Ether
400 U	UG/KG	bis(2-Chloroisopropyl) [	ther		1000 U	UG/KG	4-Nitroaniline
400 U	UG/KG	Acetophenone			1000 U	UG/KG	2-Methyl-4,6-Dinitrophenol
400 U	UG/KG	(3-and/or 4-)Methylpher		•	400 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
400 U	UG/KG	n-Nitrosodi-n-Propylam	ne		NA	UG/KG	1,2,4,5-Tetrachlorobenzene
400 U	UG/KG	Hexachloroethane			400 U	UG/KG	4-Bromophenyl Phenyl Ether
400 U	UG/KG	Nitrobenzene	·		400 U	UG/KG	Hexachlorobenzene (HCB)
400 U	UG/KG	Isophorone	*		400 U	UG/KG	Atrazine
400 U	UG/KG	2-Nitrophenol			1000 U	UG/KG	Pentachlorophenol
400 U	UG/KG	2,4-Dimethylphenol			400 U	UG/KG	Phenanthrene
400 U _	UG/KG	bis(2-Chloroethoxy)Met	nane		400 U	UG/KG	Anthracene
400 U	UG/KG	2,4-Dichlorophenol			400 U	UG/KG	Carbazole
400 U	UG/KG	Naphthalene			400 U	UG/KG	Di-n-Butylphthalate
400 U	UG/KG	4-Chloroaniline			400° U	UG/KG	Fluoranthene
400 U	UG/KG	Hexachlorobutadiene			400 U	UG/KG	Pyrene
400 U	UG/KG	Caprolactam			400 UJ	UG/KG	Benzyl Butyl Phthalate
400 U	UG/KG	4-Chloro-3-Methylpheno	ol .		400 UJ	UG/KG	3,3'-Dichlorobenzidine
400 U	UG/KG	2-Methylnaphthalene			400 U	UG/KG	Benzo(a)Anthracene
	UG/KG	Hexachlorocyclopentadi	ene (HCCP)		400 U	UG/KG	Chrysene
	UG/KG	2,4,6-Trichlorophenol			400 U	UG/KG	bis(2-Ethylhexyl) Phthalate
	UG/KG	2,4,5-Trichlorophenol			400 U	UG/KG	Di-n-Octylphthalate
	UG/KG	1,1-Biphenyl			400 U	UG/KG	Benzo(b)Fluoranthene
	UG/KG	2-Chloronaphthalene			400 U	UG/KG	Benzo(k)Fluoranthene
	UG/KG	2-Nitroaniline			400 U	UG/KG	Benzo-a-Pyrene
	UG/KG	Dimethyl Phthalate		•	400 U	UG/KG	Indeno (1,2,3-cd) Pyrene
	UG/KG	2,6-Dinitrotoluene			400 U	UG/KG	Dibenzo(a,h)Anthracene
	UG/KG	Acenaphthylene	•			.UG/KG	Benzo(ghi)Perylene
	UG/KG	3-Nitroaniline	1		18	%	% Moisture
	UG/KG-	Acenaphthene	•	•	•	•	
	UG/KG	2,4-Dinitrophenol					
1000 U	UG/KG	4-Nitrophenol					
			•				

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

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**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6763 FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS20SB /

Media: SUBSURFACE SOIL (> 12")

MD No: 1XS3

D No: 1XS3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:55

Ending:

**RESULTS UNITS ANALYTE** 

130 J UG/KG **UNKNOWN COMPOUND** 

120 NJ UG/KG **VALENCENE** 

Data Reported as Identified by CLP Lab - IDs Not Verified

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L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6764 FY 2003 Project: 03-0474

**Extractables Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS20SD / Media: SEDIMENT

Flowood, MS

Case No: 31635

MD No: 1XS4

D No: 1XS4

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:45

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
410 U	UG/KG	Benzaldehyde	410 U	UG/KG	Dibenzofuran
410 U	UG/KG	Phenol	410 U	UG/KG	2,4-Dinitrotoluene
410 U	UG/KG	bis(2-Chloroethyl) Ether	410 U	UG/KG	Diethyl Phthalate
410 U	UG/KG	2-Chlorophenol	410 U	UG/KG	Fluorene
410 U	UG/KG	2-Methylphenol	410 U	UG/KG	4-Chlorophenyl Phenyl Ether
410 U	UG/KG	bis(2-Chloroisopropyl) Ether	1000 U	UG/KG	4-Nitroaniline
410 U	UG/KG	Acetophenone	1000 U	UG/KG	2-Methyl-4,6-Dinitrophenol
410 U	UG/KG	(3-and/or 4-)Methylphenol	410 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
410 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
410 U .	UG/KG	Hexachloroethane	410 U	UG/KG	4-Bromophenyl Phenyl Ether
410 U	UG/KG	Nitrobenzene	410 U	UG/KG	Hexachlorobenzene (HCB)
410 U	UG/KG	Isophorone	410 U	UG/KG	Atrazine
410 U	UG/KG	2-Nitrophenol	1000 U	UG/KG	Pentachlorophenol
410 U	UG/KG	2,4-Dimethylphenol	410 U	UG/KG	Phenanthrene
410 U	UG/KG	bis(2-Chloroethoxy)Methane	410 U	UG/KG	Anthracene
410 U	UG/KG	2,4-Dichlorophenol	410 U	UG/KG	Carbazole
410 U	UG/KG	Naphthalene	410 U	UG/KG	Di-n-Butylphthalate
410 U	UG/KG	4-Chloroaniline	81 J	UG/KG	Fluoranthene
410 U	UG/KG	Hexachlorobutadiene	63 J	UG/KG	Pyrene
410 U	UG/KG	Caprolactam	410 UJ	UG/KG	Benzyl Butyl Phthalate
410 U	UG/KG	4-Chloro-3-Methylphenol	410 UJ	UG/KG	3,3'-Dichlorobenzidine
410 U	UG/KG	2-Methylnaphthalene	44 J	UG/KG	Benzo(a)Anthracene
410 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	69 J	UG/KG	Chrysene -
410 U	UG/KG	2,4,6-Trichlorophenol	410 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1000 U •	UG/KG	2,4,5-Trichlorophenol	410 U	UG/KG	Di-n-Octylphthalate
410 U	UG/KG	1,1-Biphenyl	140 J	UG/KG	Benzo(b)Fluoranthene
410 U	UG/KG	2-Chloronaphthalene	. 150 J	UG/KG	Benzo(k)Fluoranthene
1000 U	UG/KG	2-Nitroaniline	47 J	UG/KG	Benzo-a-Pyrene
410 U	UG/KG	Dimethyl Phthalate	51 J	UG/KG	Indeno (1,2,3-cd) Pyrene
410 U	UG/KG	2,6-Dinitrotoluene	410 U	UG/KG	Dibenzo(a,h)Anthracene
410 U	UG/KG	Acenaphthylene	410 U	UG/KG	Benzo(ghi)Perylene
1000 U	UG/KG	3-Nitroaniline	. 20	%	% Moisture
	UG/KG	Acenaphthene		•	
	UG/KG	2,4-Dinitrophenol			
1000 U	UG/KG	4-Nitrophenol			
•					
		,			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6764

FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS20SD / Media: SEDIMENT

Case No: 31635 MD No: 1XS4

Inorg Contractor: SENTIN Org Contractor: LIBRTY

D No: 1XS4

Ending:

Requestor:

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 10:45

RESULTS UNITS **ANALYTE** 

150 NJ UG/KG 220 NJ

HEXADECANOIC ACID

1-DOCOSENE UG/KG

320 NJ UG/KG STIGMAST-4-EN-3-ONE

17 UNKNOWN COMPOUNDS UG/KG 9100 J

Data Reported as Identified by CLP Lab - IDs Not Verified

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L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value,

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6765 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 10:30 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN ld/Station: GS03SD / MD No: 1XS5 Media: SEDIMENT Org Contractor: LIBRTY D No: 1XS5 DATA REPORTED ON DRY WEIGHT BASIS

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RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
110 J	UG/KG	Benzaldehyde	940 U	UG/KG	Dibenzofuran
940 U	UG/KG	Phenol	940 U	UG/KG	2,4-Dinitrotoluene
940 U	UG/KG	bis(2-Chloroethyl) Ether	940 U	UG/KG	Diethyl Phthalate
940 U	UG/KG	2-Chlorophenol	130 U	UG/KG	Fluorene
940 U	UG/KG	2-Methylphenol	940 U	UG/KG	4-Chlorophenyl Phenyl Ether
940 U	UG/KG	bis(2-Chloroisopropyl) Ether	2400 U	UG/KG	4-Nitroaniline
940 U ·	UG/KG	Acetophenone	2400 U	UG/KG	2-Methyl-4,6-Dinitrophenol
940 U	UG/KG	(3-and/or 4-)Methylphenol	940 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
940 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
940 U	UG/KG	Hexachloroethane	940 U	UG/KG	4-Bromophenyl Phenyl Ether
940 U	UG/KG	Nitrobenzene	940 U	UG/KG	Hexachlorobenzene (HCB)
940 U	UG/KG	Isophorone	940 U	UG/KG	Atrazine
940 U	UG/KG	2-Nitrophenol	190 U	UG/KG	Pentachlorophenol
940 U	UG/KG	2,4-Dimethylphenol	- 120 J	UG/KG	Phenanthrene
940 U	UG/KG	bis(2-Chloroethoxy)Methane	270 J	UG/KG	Anthracene
940 U	UG/KG	2,4-Dichlorophenol	940 U	UG/KG	Carbazole
940 U	UG/KG	Naphthalene	940 U	UG/KG	Di-n-Butylphthalate
. 940 U	UG/KG	4-Chloroaniline	340 J	UG/KG	Fluoranthene
940 U	UG/KG	Hexachlorobutadiene	220 J	UG/KG	Pyrene
940 U	UG/KG	Caprolactam	940 UJ	UG/KG	Benzyl Butyl Phthalate
940 U	UG/KG	4-Chloro-3-Methylphenol	940 UJ	UG/KG	3.3'-Dichlorobenzidine
930 U	UG/KG	2-Methylnaphthalene	150 J	UG/KG	Benzo(a)Anthracene
940 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	250 J	UG/KG	Chrysene
940 U	UG/KG	2,4,6-Trichlorophenol	940 U	UG/KG	bis(2-Ethylhexyl) Phthalate
2400 U	UG/KG	2,4,5-Trichlorophenol	940 U	UG/KG	Di-n-Octylphthalate
940 U	UG/KG	1,1-Biphenyl	540 J	UG/KG	Benzo(b)Fluoranthene
940 U	UG/KG	2-Chloronaphthalene	560 J	UG/KG	Benzo(k)Fluoranthene
. 2400 U	UG/KG	2-Nitroaniline	. 150 J	UG/KG	Benzo-a-Pyrene
940 U	UG/KG	Dimethyl Phthalate	180 J	UG/KG	Indeno (1,2,3-cd) Pyrene .
940 U	UG/KG	2,6-Dinitrotoluene	940 U	UG/KG	Dibenzo(a,h)Anthracene
940 U	UG/KG	Acenaphthylene	940 U	UG/KG	Benzo(ghi)Perylene
2400 U	UG/KG	3-Nitroaniline	79	%	% Moisture
940 U	UG/KG	Acenaphthene			
2400 U	UG/KG	2,4-Dinitrophenol			
2400 U	UG/KG	4-Nitrophenol		•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

6765 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 10:30

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Sample

Case No: 31635

Id/Station: GS03SD /

MD No: 1XS5

Inorg Contractor: SENTIN

Media: SEDIMENT

D No: 1XS5

Org Contractor: LIBRTY

RESULTS UNITS

590 NJ UG/KG 1000 NJ UG/KG 1400 NJ UG/KG

**ANALYTE** BENZO [E] PYRENE

1-HEXADECENE STIGMAST-4-EN-3-ONE

54000 J UG/KG 25 UNKNOWNS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

FY 2003

Project: 03-0474

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS03SD /

Case No: 31635

Media: SEDIMENT

MD No: 1XS5

D No: 1XS5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:30

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
33	UG/KG	2-Methylnaphthalene
51	UG/KG	Naphthalene
44	UG/KG	Acenaphthylene
100	UG/KG	Acenaphthene
82	UG/KG	Fluorene
72	UG/KG	Phenanthrene
140	UG/KG	Anthracene
190	UG/KG	Fluoranthene
140	UG/KG	Pyrene
110	UG/KG	Benzo(a)Anthracene
200	UG/KG	Chrysene
NA	UG/KG	Benzo(b)Fluoranthene
NA	UG/KG	Benzo(k)Fluoranthene
120	UG/KG	Benzo-a-Pyrene
NA	UG/KG	Indeno (1,2,3-cd) Pyrene
40 .	UG/KG	Dibenzo(a,h)Anthracene
NA	UG/KG	Benzo(ghi)Perylene
100 U	UG/KG	Pentachlorophenol
		<b>\</b>

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

FY 2003 Project: 03-0474 6766 Sample **Extractables Scan** 

Flowood, MS

Case No: 31635

Produced by: Goddard, Denise

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:30

Program: SF

Ending:

ld/Station: GS16SS /

Inorg Contractor: SENTIN MD No: 1XS6

D No: 1XS6

Media: SURFACE SOIL (0" - 12")

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE		•		RESULTS	UNITS	ANALYTE	
: 420 U	UG/KG	Benzaldehyde	i -			420 U	UG/KG	Dibenzofuran	
420 U	UG/KG	Phenol				420 U	UG/KG	2,4-Dinitrotoluene	
420 U	UG/KG	bis(2-Chloroethyl) Ethe	r			420 U	UG/KG	Diethyl Phthalate	
420 U	UG/KG	2-Chlorophenol	:		•	420 U	UG/KG	Fluorene	
420 U	UG/KG	2-Methylphenol	!			420 U	UG/KG	4-Chlorophenyl Phenyl Ether	
420 U	UG/KG	bis(2-Chloroisopropyl) E	ther			1100 U	UG/KG	4-Nitroaniline	
420 U	UG/KG	Acetophenone				1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol	
420 U	UG/KG	(3-and/or 4-)Methylpher				420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine	
420 U	UG/KG	n-Nitrosodi-n-Propylam	ine		в .	NA	UG/KG	1,2,4,5-Tetrachlorobenzene	
420 U	UG/KG	Hexachloroethane				420 U	UG/KG	4-Bromophenyl Phenyl Ether	
420 U	UG/KG	Nitrobenzene	:			420 U	UG/KG	Hexachlorobenzene (HCB)	
420 U	UG/KG	Isophorone	į			420 U	UG/KG	Atrazine	
420 U	UG/KG	2-Nitrophenol	! -			1100 U	UG/KG	Pentachlorophenol	
420 U	UG/KG	2,4-Dimethylphenol	•			310 J	UG/KG	Phenanthrene	
420 U	UG/KG	bis(2-Chloroethoxy)Met	hane			420 U	UG/KG	Anthracene	
420 U	UG/KG	2,4-Dichlorophenol				62 J	UG/KG	Carbazole	
420 U	UG/KG	Naphthalene	i'			420 U	UG/KG	Di-n-Butylphthalate	
420 U	UG/KG	4-Chloroaniline				760	UG/KG	Fluoranthene	
420 U	UG/KG	Hexachlorobutadiene		•		540	UG/KG	Pyrene	
420 U	UG/KG	Caprolactam		•		420 UJ	UG/KG	Benzyl Butyl Phthalate	
420 U	UG/KG	4-Chloro-3-Methylpheno	ol .	•		420 UJ	UG/KG	3,3'-Dichlorobenzidine	
420 U	UG/KG	2-Methylnaphthalene	i i			260 J	UG/KG	Benzo(a)Anthracene	
420 U	UG/KG	Hexachlorocyclopentadi	ene (HCCP)			390 J	UG/KG	Chrysene	
420 U	UG/KG	2,4,6-Trichlorophenol	!			420 U	UG/KG	bis(2-Ethylhexyl) Phthalate	•
1100 U	UG/KG	2,4,5-Trichlorophenol				420 U	UG/KG	Di-n-Octylphthalate	
420 U	UG/KG	1,1-Biphenyl				700 J	UG/KG	Benzo(b)Fluoranthene	
420 U	UG/KG	2-Chloronaphthalene		•	-	. 730 J	UG/KG	Benzo(k)Fluoranthene	
1100 U	UG/KG	2-Nitroaniline				230 J	UG/KG	Benzo-a-Pyrene	
420 U	UG/KG	Dimethyl Phthalate	i			270 J	UG/KG	Indeno (1,2,3-cd) Pyrene	
420 U	UG/KG	2,6-Dinitrotoluene				95 J	UG/KG	Dibenzo(a,h)Anthracene	
	UG/KG	Acenaphthylene	1			92 J	UG/KG	Benzo(ghi)Perylene	
	UG/KG	3-Nitroaniline	i			21	%	% Moisture	
	UG/KG	Acenaphthene							
1100 U	UG/KG	2,4-Dinitrophenol	!						
1100 U ·	UG/KG	4-Nitrophenol							•
			1						

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6766 FY 2003

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS16SS /

MD No: 1XS6

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XS6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:30

Ending:

RESULTS	UNITS	ANALYTE
160 J	UG/KG	9,10-ANTHRACENEDIONE
2000 J	UG/KG	PYRENE, 1-METHYL-
1400 J	UG/KG	PYRENE, 2-METHYL-
1400 J	UG/KG	BENZO [B] NAPHTHO [2,1-D] THIOPH
1600 J	UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
2300 J	UG/KG	CHRYSENE, 5-METHYL-
2600 J	UG/KG	5,12-NAPHTHACENEDIONE
.200 J	UG/KG	PERYLENE
580 J	UG/KG	UNKNOWN PAH
1200 NJ	UG/KG	VALENCENE
24000 J	`UG/KG	17 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Beporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample FY 2003 6767

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS16SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

MD No: 1XS7

D No: 1XS7

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

DECLU TO	LINUTO	ANALVE	DECLU TO	LIMITO	ANAL VTF
RESULTS	UNITS	ANALYTE	RESULTS		ANALYTE
420 U	UG/KG	Benzaldehyde	420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	420 U	UG/KG	Diethyl Phthalate
420 U	UG/KG	2-Chlorophenol	420 U	UG/KG	Fluorene
420 U	UG/KG	2-Methylphenol	420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
420 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
420 U	UG/KG	(3-and/or 4-)Methylphenol	420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane	420 U	UG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	420 U	UG/KG	Hexachlorobenzene (HCB)
420 U	UG/KG	Isophorone	420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
420 U	UG/KG	2,4-Dimethylphenol	420 U	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Methane	420 U	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol	420 U	UG/KG	Carbazole
420 U.	UG/KG	Naphthalene	420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline	420 U	UG/KG	Fluoranthene
420 U	UG/KG	Hexachlorobutadiene*	420 U	UG/KG	Pyrene
420 U	UG/KG	Caprolactam	420 UJ	UG/KG	Benzyl Butyl Phthalate
420 U	UG/KG	4-Chloro-3-Methylphenol	420 UJ	UG/KG	3.3'-Dichlorobenzidine
420 U	UG/KG	2-Methylnaphthalene	420 U	UG/KG	Benzo(a)Anthracene
420 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	420 U	UG/KG	Chrysene
420 U	UG/KG	2,4,6-Trichlorophenol	420 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	420 U	UG/KG	Di-n-Octylphthalate
420 U	UG/KG	1,1-Biphenyl	420 U .	UG/KG	Benzo(b)Fluoranthene
420 U	UG/KG	2-Chloronaphthalene	420 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	420 U	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate	420 <sup>'</sup> U	UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2,6-Dinitrotoluene	420 U	UG/KG	Dibenzo(a,h)Anthracene
420 U	UG/KG	Acenaphthylene	420 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	22	%	% Moisture
420 U	UG/KG	Acenaphthene	— <del></del>		,
1100 U	UG/KG	2,4-Dinitrophenol			•
1100 U	UG/KG	4-Nitrophenol			
	-				·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates...

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

EXTRACTABLES SAMPLE ANALYSIS

Production Date: 06/20/2003 14:10

Sample

6767 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS16SB / MD No: 1XS7

Media: SUBSURFACE SOIL (> 12") D No: 1XS7 Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:40

Ending:

Requestor:

**RESULTS UNITS** 

220 J

**ANALYTE** UG/KG

2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

6768 FY 2003 Project: 03-0474 Sample **Extractables Scan** Flowood, MS Produced by: Goddard, Denise Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:05

Ending:

Id/Station: GS21SS /

Program: SF

Case No: 31635

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

Facility: Gulf States Creosoting

MD No: 1XS8 Org Contractor: LIBRTY D No: 1XS8

DATA REPORTED ON DRY WEIGHT BASIS

<del></del>	<del></del>	,			
RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
420 U	UG/KG	Benzaldehyde	420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	420 U	UG/KG	Diethyl Phthalate
420 U	UG/KG	2-Chlorophenol	420 U	UG/KG	Fluorene
420 U	UG/KG	2-Methylphenol	420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
420 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
420 U	UG/KG	(3-and/or 4-)Methylphenol	420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane	420 U	UG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	420 U	UG/KG	Hexachlorobenzene (HCB)
420 U	UG/KG	Isophorone	420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
420 U	UG/KG	2,4-Dimethylphenol	420 U	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Methane	420 U	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol	420 U	UG/KG	Carbazole
420 U	UG/KG	Naphthalene	420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline	420 U	UG/KG	Fluoranthene
420 U	UG/KG	Hexachlorobutadiene	420 U	UG/KG	Pyrene
420 U	UG/KG	Caprolactam	420 UJ	UG/KG	Benzyl Butyl Phthalate
420 U	UG/KG	4-Chloro-3-Methylphenol	420 UJ	UG/KG	3.3'-Dichlorobenzidine
420 U	UG/KG	2-Methylnaphthalene	420 U	UG/KG	Benzo(a)Anthracene
420 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	420 U	UG/KG	Chrysene
420 U	UG/KG	2,4,6-Trichlorophenol	420 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	420 U	UG/KG	Di-n-Octylphthalate
420 U	UG/KG	1,1-Biphenyl	420 U	UG/KG	Benzo(b)Fluoranthene
420 U*	UG/KG	2-Chloronaphthalene	420 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	420 U	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate	420 U	UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2.6-Dinitrotoluene	420 U	UG/KG	Dibenzo(a,h)Anthracene
420 U	UG/KG	Acenaphthylene	420 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	22	%	% Moisture
420 U	UG/KG	Acenaphthene			, c 1110101010
1100 U	UG/KG	2,4-Dinitrophenol			
1100 U	UG/KG	4-Nitrophenol			
			,		
				•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6768 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS21SS /

MD No: 1XS8 D No: 1XS8

Media: SURFACE SOIL (0" - 12")

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:05

Ending:

**RESULTS UNITS** 230 NJ UG/KG

**ANALYTE** 

1-EICOSANOL

3100 J UG/KG 14 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Project: 03-0474 Sample 6769 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 12:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN ld/Station: GS21SB / MD No: 1XS9 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") **D No: 1XS9** DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
410 U	UG/KG	Benzaldehyde	410 U	UG/KG	Dibenzofuran
410 U	UG/KG	Phenol	410 U	UG/KG	2,4-Dinitrotoluene
410 U	UG/KG	bis(2-Chloroethyl) Ether	410 U	UG/KG	Diethyl Phthalate
410 U	UG/KG	2-Chlorophenol	410 U	UG/KG	Fluorene
410 U	UG/KG	2-Methylphenol	410 U	UG/KG	4-Chlorophenyl Phenyl Ether
410 U	UG/KG	bis(2-Chloroisopropyl) Ether	1000 U	UG/KG	4-Nitroaniline
410 U	UG/KG	Acetophenone	1000 U	UG/KG	2-Methyl-4,6-Dinitrophenol
410 U	UG/KG	(3-and/or 4-)Methylphenol	410 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
410 U *	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
410 U	UG/KG	Hexachloroethane	410 U	UG/KG	4-Bromophenyl Phenyl Ether
410 U	UG/KG	Nitrobenzene	410 U	UG/KG	Hexachlorobenzene (HCB)
. 410 U	UG/KG	Isophorone	410 U	UG/KG	Atrazine
410 U	UG/KG	2-Nitrophenol	1000 U	UG/KG	Pentachlorophenol
410 U	UG/KG	2,4-Dimethylphenol	410 U	UG/KG	Phenanthrene
410 U	UG/KG	bis(2-Chloroethoxy)Methane	410 U	UG/KG	Anthracene
410 U	UG/KG	2,4-Dichlorophenol	410 U	UG/KG	Carbazole
410 U	UG/KG	Naphthalene	410 U	UG/KG	Di-n-Butylphthalate
410 U	UG/KG	4-Chloroaniline	410 U	UG/KG	Fluoranthene
410 U	UG/KG	Hexachlorobutadiene	410 U	UG/KG	Pyrene
410 U	UG/KG	Caprolactam	410 UJ.	UG/KG	Benzyl Butyl Phthalate
410 U	UG/KG	4-Chloro-3-Methylphenol	410 UJ	UG/KG	3,3'-Dichlorobenzidine
410 U	UG/KG	2-Methylnaphthalene	410 U	UG/KG	Benzo(a)Anthracene
410 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	410 U	UG/KG	Chrysene
410 U	UG/KG	2,4,6-Trichlorophenol	410 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1000 U	UG/KG	2,4,5-Trichlorophenol	410 U	UG/KG	Di-n-Octylphthalate
410 U	UG/KG	1,1-Biphenyl	410 U	UG/KG	Benzo(b)Fluoranthene
410 U	UG/KG	2-Chloronaphthalene	410 U	UG/KG	Benzo(k)Fluoranthene
1000 U	UG/KG	2-Nitroaniline	410 U	UG/KG	Benzo-a-Pyrene
410 U	UG/KG	Dimethyl Phthalate	410 U	UG/KG	Indeno (1,2,3-cd) Pyrene
410 U	UG/KG	2,6-Dinitrotoluene	410 U	UG/KG	Dibenzo(a,h)Anthracene
410 U	UG/KG	Acenaphthylene	410 U	UG/KG	Benzo(ghi)Perylene
1000 U	UG/KG	3-Nitroaniline	20	%	% Moisture
410 U	UG/KG	Acenaphthene			
1000 U	UG/KG	2,4-Dinitrophenol			·
1000 U	UG/KG	4-Nitrophenol			• •
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•		Y .			·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 12:15

Requestor:

Ending:

Sample

6769 FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS21SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

D No: 1XS9

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

MD No. 1XS9

RESULTS UNITS **ANALYTE** 

> 84 J UG/KG

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value,

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample\* 6770 FY 2003

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS01GW /

MD No: 1XT0

Media: GROUNDWATER

Case No: 31635

D No: 1XT0

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:05

Ending:

RESULTS	UNITS	ANALYTE			RESULTS		ANALYTE	
5.0 UJ	UG/L	Benzaldehyde			5.0 UJ	UG/L	Dibenzofuran	
5.0 UJ	UG/L	Phenol			5.0 UJ	UG/L	2,4-Dinitrotoluene	•
5.0 UJ	UG/L	bis(2-Chloroethyl) Ether			5.0 UJ	UG/L	Diethyl Phthalate	
5.0 UJ	UG/L	2-Chlorophenol			5.0 UJ	UG/L	Fluorene	
5.0 UJ	UG/L .	2-Methylphenol			5.0 UJ	UG/L	4-Chlorophenyl Phenyl Ether	
5.0 UJ	UG/L	bis(2-Chloroisopropyl) Eth	er .		20 UJ	UG/L	4-Nitroaniline	
5.0 UJ	UG/L	Acetophenone			20 UJ	UG/L	2-Methyl-4,6-Dinitrophenol	
5.0 UJ	UG/L	(3-and/or 4-)Methylphenol			5.0 UJ	UG/L	n-Nitrosodiphenylamine/Diphenylamine	
5.0 UJ	UG/L	n-Nitrosodi-n-Propylamine		-	5.0 UJ	UG/L	1,2,4,5-Tetrachlorobenzene	
5.0 UJ	UG/L	Hexachloroethane			5.0 UJ	UG/L	4-Bromophenyl Phenyl Ether	
5.0 UJ	UG/L	Nitrobenzene			5.0 UJ	UG/L	Hexachlorobenzene (HCB)	
5.0 UJ	UG/L	Isophorone			5.0 UJ	UG/L	Atrazine	
5.0 UJ	UG/L	2-Nitrophenol	•		5.0 UJ	UG/L	Pentachlorophenol	•
5.0 UJ	UG/L	2,4-Dimethylphenol	·		5.0 UJ	UG/L	Phenanthrene	•
5.0 UJ	UG/L	bis(2-Chloroethoxy)Metha	ne		5.0 UJ	UG/L	Anthracene	
5.0 UJ	UG/L	2,4-Dichlorophenol	•		NA	·UG/L	Carbazole	
5.0 UJ	UG/L	Naphthalene	•	•	5.0 UJ	UG/L	Di-n-Butylphthalate	
5.0 UJ	UG/L	4-Chloroaniline		•	5.0 UJ	UG/L	Fluoranthene	
5.0 UJ	UG/L	Hexachlorobutadiene			5.0 UJ	UG/L	Pyrene	
5.0 UJ	UG/L	Caprolactam			5.0 UJ	UG/L	Benzyl Butyl Phthalate	
5.0 UJ	UG/L	4-Chloro-3-Methylphenol			5.0 UJ	UG/L	3,3'-Dichlorobenzidine	
5.0 UJ	UG/L	2-Methylnaphthalene			5.0 UJ	UG/L	Benzo(a)Anthracene	
5.0 UJ	UG/L	Hexachlorocyclopentadien	e (HCCP)		5.0 UJ	UG/L	Chrysene	
5.0 UJ	UG/L	2,4,6-Trichlorophenol			5.0 UJ	UG/L	bis(2-Ethylhexyl) Phthalate	
20 ÚJ	UG/L	2,4,5-Trichlorophenol			5.0 UJ	UG/L	Di-n-Octylphthalate	
5.0 UJ	UG/L	1,1-Biphenyl		•	5.0 UJ	UG/L	Benzo(b)Fluoranthene	•
5.0 UJ	UG/L	2-Chloronaphthalene			5.0 UJ	UG/L	Benzo(k)Fluoranthene	
20 UJ	UG/L	2-Nitroaniline			5.0 UJ	UG/L	Benzo-a-Pyrene	
5.0 UJ	UG/L	Dimethyl Phthalate			5.0 UJ	UG/L	Indeno (1,2,3-cd) Pyrene	
5.0 UJ	UG/L	2,6-Dinitrotoluene			5.0 UJ	UG/L	Dibenzo(a,h)Anthracene	
5.0 UJ	UG/L	Acenaphthylene			5.0 UJ	UG/L	Benzo(ghi)Perylene	
20 UJ	UG/L	3-Nitroaniline						
5.0 UJ	UG/L	Acenaphthene			·		•	
20 UJ	UG/L	2,4-Dinitrophenol	•					
20 UJ	UG/L	4-Nitrophenol		•				
•		į					•	
							•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6770 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS01GW /

Media: GROUNDWATER

Flowood, MS

Case No: 31635 MD No: 1XT0

D No: 1XT0

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:05

Ending:

**RESULTS UNITS ANALYTE** 2.2 NJ UG/L **D-LIMONENE** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6771 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS01SB / MD No: 1XT1 Media: SUBSURFACE SOIL (> 12") Org Contractor: LIBRTY D No: 1XT1 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
400 U	UG/KG	Benzaldehyde	400 U	UG/KG	Dibenzofuran
400 U	UG/KG <sup>,</sup>	Phenol	400 U	UG/KG	2,4-Dinitrotoluene
400 U	UG/KG	bis(2-Chloroethyl) Ether	400 U	UG/KG	Diethyl Phthalate
400 U	UG/KG	2-Chlorophenol	400 U	UG/KG	Fluorene
400 U	UG/KG	2-Methylphenol	400 U	UG/KG	4-Chlorophenyl Phenyl Ether
400 U	UG/KG	bis(2-Chloroisopropyl) Ether	1000 U	UG/KG	4-Nitroaniline
400 U	UG/KG	Acetophenone	1000 U	UG/KG	2-Methyl-4,6-Dinitrophenol
400 U	UG/KG	(3-and/or 4-)Methylphenol	400 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
400 U	UG/KG	n-Nitrosodi-n-Propylamine	. NA	UG/KG	1,2,4,5-Tetrachlorobenzene
400 U	UG/KG	Hexachloroethane	400 U	UG/KG	4-Bromophenyl Phenyl Ether
400 U	UG/KG	Nitrobenzene	400 U	UG/KG	Hexachlorobenzene (HCB)
400 U	UG/KG	Isophorone	400 U	UG/KG	Atrazine
400 U	UG/KG	2-Nitrophenol	1000 U	UG/KG	Pentachlorophenol
400 U	UG/KG	2,4-Dimethylphenol	400 U	UG/KG	Phenanthrene
400 U	UG/KG	bis(2-Chloroethoxy)Methane	400 U	UG/KG	Anthracene
400 U	UG/KG	2,4-Dichlorophenol	400 U	UG/KG	Carbazole
400 U	UG/KG	Naphthalene	400 U	UG/KG	Di-n-Butylphthalate
400 U	UG/KG	4-Chloroaniline	400 U	UG/KG	Fluoranthene
400 U	UG/KG	Hexachlorobutadiene	400 U	UG/KG	Pyrene
400 U	UG/KG	Caprolactam	400 UJ	UG/KG	Benzyl Butyl Phthalate
400 U	UG/KG	4-Chloro-3-Methylphenol	400 UJ	UG/KG	3,3'-Dichlorobenzidine
400 U	UG/KG	2-Methylnaphthalene	400 U	UG/KG	Benzo(a)Anthracene
400 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	400 U	UG/KG	Chrysene
400 U	UG/KG	2,4,6-Trichlorophenol	400 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1000 U	UG/KG	2,4,5-Trichlorophenol	400 U	UG/KG	Di-n-Octylphthalate
400 U	UG/KG	1,1-Biphenyl	400 U	UG/KG	Benzo(b)Fluoranthene
400 U	UG/KG	2-Chloronaphthalene	400 U	UG/KG	Benzo(k)Fluoranthene
1000 U *	UG/KG	2-Nitroaniline	400 U	UG/KG	Benzo-a-Pyrene
· 400 U	UG/KG	Dimethyl Phthalate	400 U	UG/KG	Indeno (1,2,3-cd) Pyrene
400 U	UG/KG	2,6-Dinitrotoluene	400 U	UG/KG	Dibenzo(a,h)Anthracene
400 U	UG/KG	Acenaphthylene	400 U	UG/KG	Benzo(ghi)Perylene
1000 U	UG/KG	3-Nitroaniline	18	%	% Moisture
400 U	UG/KG	Acenaphthene			·
1000 U	UG/KG	2,4-Dinitrophenol			
1000 U	UG/KG	4-Nitrophenol			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6771 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 09:15

Ending:

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

MD No: 1XT1

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

D No: 1XT1

Org Contractor: LIBRTY

**RESULTS UNITS** 

92 NJ

Id/Station: GS01SB /

**ANALYTE** UG/KG

1-PHENANTHRENECARBOXYLIC ACID

UG/KG **5 UNKNOWN COMPOUNDS** 840 J

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6772 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:00 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS01SS / MD No: 1XT2 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT2 DATA REPORTED ON DRY WEIGHT BASIS

380 U   UG/KG   Benzaldehyde   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrolubuene   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol   380 U   UG/KG   A-Dinitrophenol	RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
380 U   UG/KG   UG/KG   UG/KG   C-Chloropethyl) Ether   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-Chlorophenol   380 U   UG/KG   C-CCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCCC	380 U	UG/KG	Benzaldehyde	380 U	UG/KG	
380 U   UG/KG   2-Methylphenol   380 U   UG/KG   4-Chlorophenyl Phenyl Ether   380 U   UG/KG   4-Chlorophenyl Phenyl Ether   380 U   UG/KG   4-Chlorophenyl Phenyl Ether   380 U   UG/KG   4-Nitroaniline   380 U   UG/KG   4-Nitroaniline   380 U   UG/KG   4-Nitroaniline   380 U   UG/KG   3-Methylphenol   380 U   UG/KG   3-Methylphenol   380 U   UG/KG   1-Ritrosodi-n-Propylamine   12,4,5-Tetrachlorobenzene   12,4,5-Tetrachlorobenzene   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritrosodi-n-Propylamine   380 U   UG/KG   1-Ritros		UG/KG	Phenol	U 086	UG/KG	2,4-Dinitrotoluene
380 U   UG/KG   2-Methylphenol   380 U   UG/KG   4-Chlorophenyl Phenyl Ether   380 U   UG/KG   4-Chlorophenyl Phenyl Ether   380 U   UG/KG   4-Chlorophenyl Phenyl Ether   380 U   UG/KG   4-Mitroaniline   2-Methyl-4,6-Dinitrophenol   380 U   UG/KG   4-Mitroaniline   2-Methyl-4,6-Dinitrophenol   380 U   UG/KG   3-and/or 4-Methylphenol   380 U   UG/KG   1-Nitrosodi-n-Propylamine   1,2,4,5-Tetrachiorobenzene   380 U   UG/KG   1-Nitrosodi-n-Propylamine   1,2,4,5-Tetrachiorobenzene   380 U   UG/KG   1-Nitrosodi-n-Propylamine   1,2,4,5-Tetrachiorobenzene   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propylamine   380 U   UG/KG   1-Nitrosodi-n-Propyl	380 U	UG/KG	bis(2-Chloroethyl) Ether	380 U	UG/KG	Diethyl Phthalate
380 U   UG/KG   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary   Carbonary		UG/KG		380 U	UG/KG	Fluorene
380 U   UG/KG   Dis(2-Chloroisopropyl) Ether   970 U   UG/KG   Chloroisopropyl) henyl Ether   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4,5-Tetrachlorobenzene   1,2,4		UG/KG	2-Methylphenol	380 U	UG/KG	4-Chlorophenyl Phenyl Ether
380 U         UG/KG         Acetophenone         970 U         UG/KG         2-Methyl-4,6-Dinitrophenol           380 U         UG/KG         1,000         1,000         1,000         1,2,4,5-Tetrachlorobenzene           380 U         UG/KG         1,2,4,5-Tetrachlorobenzene         4-Bromophenyl Phenyl Ether           380 U         UG/KG         1,2,4,5-Tetrachlorobenzene         4-Bromophenyl Phenyl Ether           380 U         UG/KG         1,2,4,5-Tetrachlorobenzene         4-Bromophenyl Phenyl Ether           380 U         UG/KG         1,000         4-Bromophenyl Phenyl Ether           4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether           380 U         UG/KG         2,4-Dintrophenol         380 U         UG/KG         Atrazine           380 U         UG/KG         2,4-Dintrophenol         380 U         UG/KG         Pentachlorophenol           380 U         UG/KG         2,4-Dichlorophenol         380 U         UG/KG         Carbazole           380 U         UG/KG         Naphthalene         380 U         UG/KG         Carbazole           380 U         UG/KG         AChloropathilalene         380 U         UG/KG         Benzyl Bullyl Phithalate           380 U         UG/KG		UG/KG	bis(2-Chloroisopropyl) Ether	970 UJ	UG/KG	4-Nitroaniline
380 U         UG/KG 380 U         (3-and/or 4-)Methylphenol         380 U         UG/KG NA         n-hitrosodiphenylamine/Diphenylamine         1,2,4,5-Tetrachlorobenzene         1,2,4,5-Tetrachlorobenzene         1,2,4,5-Tetrachlorobenzene         1,2,4,5-Tetrachlorobenzene         1,2,4,5-Tetrachlorobenzene         1,2,4,5-Tetrachlorobenzene         1,2,4,5-Tetrachlorobenzene         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl Phenyl Ether         4-Bromophenyl		UG/KG		970 U	UG/KG	2-Methyl-4,6-Dinitrophenol
380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Hexachlorocethane   380 U   UG/KG   Pentachlorophenol   380 U   UG/KG   Pentachlorophenol   380 U   UG/KG   Pentachlorophenol   380 U   UG/KG   Anthracene   380 U   UG/KG   Anthracene   380 U   UG/KG   Anthracene   380 U   UG/KG   Anthracene   380 U   UG/KG   Hexachlorobutadiene   380 U   UG/KG   Hexachlorobutadiene   380 U   UG/KG   Hexachlorobutadiene   380 U   UG/KG   Pyrene   380 U   UG/KG   Hexachlorocothadiene   380 U   UG/KG   Pyrene   380 U   UG/KG   Hexachlorocothadiene   380 U   UG/KG   San Dichlorobenzidine   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene   380 U   UG/KG   San Dichlorocothadiene		UG/KG <sup>*</sup>	(3-and/or 4-)Methylphenol	380 U	UG/KG	
380 U         UG/KG Johnsone         380 UJ UG/KG Johnsone         Hexachlorobenzene (HCB)           380 U         UG/KG Johnsone         380 UJ UG/KG Johnsone         Hexachlorophenol           380 U         UG/KG Johnsone         2-Nitrophenol         380 UJ UG/KG Pentachlorophenol           380 U         UG/KG Johnsone         2-4-Dimethylphenol         380 UJ UG/KG Phenanthrene           380 U         UG/KG Johnsone         380 UJ UG/KG Phenanthrene           380 U         UG/KG Johnsone         380 UJ UG/KG Phenanthrene           380 U         UG/KG UG/KG Johnsone         2-4-Dichlorophenol         380 UJ UG/KG Carbazole           380 U         UG/KG Johnsone         380 UJ UG/KG UG/KG Din-Butylphthalate         380 UJ UG/KG Din-Butylphthalate           380 U         UG/KG Johnsone         380 UJ UG/KG UG/KG Din-Butylphthalate         380 UJ UG/KG Pyrene           380 U         UG/KG Johnsone         380 UJ UG/KG Din-Butylphthalate         380 UJ UG/KG Benzyl Butyl Phthalate           380 UJ UG/KG Johnsone         2-Methylnaphthalene         380 UJ UG/KG Benzyl Butyl Phthalate         380 UJ UG/KG Benzyl Butyl Phthalate           380 UJ UG/KG Johnsone         2-4,5-Trichlorophenol         380 UJ UG/KG Benzyl Butyl Phthalate         380 UJ UG/KG Benzyl Butyl Phthalate           380 UJ UG/KG Johnsone         2-4,5-Trichlorophenol         380 UJ UG/KG Benzyl Buty		UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
380 U         UG/KG         Isophorone         380 U         UG/KG         Atrazine           380 U         UG/KG         2-Nitrophenol         380 U         UG/KG         Pentachlorophenol           380 U         UG/KG         2,4-Dimethylphenol         380 U         UG/KG         Anthracene           380 U         UG/KG         2,4-Dichlorophenol         380 U         UG/KG         Anthracene           380 U         UG/KG         Naphthalene         380 U         UG/KG         Carbazole           380 U         UG/KG         Hexachlorobutadiene         380 U         UG/KG         Fluoranthene           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         4-Chloro-3-Methylphenol         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         4-Chloro-a-Methylphenol         380 U         UG/KG         Discoplanting           380 U         UG/KG         2,4,5-Trichlorophenol <td>380 U*</td> <td>UG/KG</td> <td>Hexachloroethane</td> <td>380 U</td> <td>UG/KG</td> <td>4-Bromophenyl Phenyl Ether</td>	380 U*	UG/KG	Hexachloroethane	380 U	UG/KG	4-Bromophenyl Phenyl Ether
380 U         UG/KG         2-Nitrophenol         970 UJ         UG/KG         Pentachlorophenol           380 U         UG/KG         2,4-Dimethylphenol         380 U         UG/KG         Phenanthrene           380 U         UG/KG         2,4-Dichlorophenol         380 U         UG/KG         Anthracene           380 U         UG/KG         2,4-Dichlorophenol         380 U         UG/KG         Carbazole           380 U         UG/KG         4-Chloroaniline         380 U         UG/KG         Din-Butylphthalate           380 U         UG/KG         4-Chloroaniline         380 U         UG/KG         Fluoranthene           380 U         UG/KG         Hexachlorobutadiene         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         4-Chloro-3-Methylphenol         380 U         UG/KG         3-3-Dichlorobenzidine           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-A,6-Trichlorophenol         380 U         UG/KG         Benzo(a)Anthracene           970 U         UG/KG         2-A,5-Trichlorophenol         380 U         UG/KG         Benzo(b)Fluoranthene           380 U         UG/KG	380 U	UG/KG	Nitrobenzene		UG/KG	Hexachlorobenzene (HCB)
380 U         UG/KG         2.4-Dimethylphenol         380 U         UG/KG         Anthracene           380 U         UG/KG         2.4-Dichloropthoxy)Methane         380 U         UG/KG         Anthracene           380 U         UG/KG         2.4-Dichlorophenol         380 U         UG/KG         Carbazole           380 U         UG/KG         Naphthalene         380 U         UG/KG         Di-n-Butylphthalate           380 U         UG/KG         4-Chloroanline         380 U         UG/KG         Fluoranthene           380 U         UG/KG         Hexachlorobutatiene         380 U         UG/KG         Pyrene           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-4,6-Trichlorophenol         380 U         UG/KG         Chrysene           380 U         UG/KG         2,4-5-Trichlorophenol         380 U         UG/KG         Di-n-Octylphthalate           380 U         UG/KG         2,5-Trichlor		UG/KG	Isophorone	380 U	UG/KG	Atrazine
380 U         UG/KG         bis(2-Chloroethoxy)Methane         380 U         UG/KG         Anthracene           380 U         UG/KG         2,4-Dichlorophenol         380 U         UG/KG         Carbazole           380 U         UG/KG         Naphthalene         380 U         UG/KG         Din-Butylphthalate           380 U         UG/KG         4-Chloroaniline         380 U         UG/KG         Fluoranthene           380 U         UG/KG         Aprolactam         380 U         UG/KG         Pyrene           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Chrysene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Dis(2-Ethylhexyl) Phthalate           970 U         UG/KG         2-Methylnaph	380 U	UG/KG	2-Nitrophenol	- 970 UJ	UG/KG	Pentachlorophenol
380 U         UG/KG         2,4-Dichlorophenol         380 U         UG/KG         Carbazole           380 U         UG/KG         Naphthalene         380 U         UG/KG         Di-n-Butylphthalate           380 U         UG/KG         4-Chloroaniline         380 U         UG/KG         Fluoranthene           380 U         UG/KG         Hexachlorobutadiene         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Kethylnaphthalene         380 U         UG/KG         Chrysene           380 U         UG/KG         2-Kethylnaphthalene         380 U         UG/KG         Di-n-Octylphthalate           380 U         UG/KG         2-4,6-Trichlorophenol         380 U         UG/KG         Di-n-Octylphthalate           970 U         UG/KG         2-4,5-Trichlorophenol         380 U         UG/KG         Benzo(b)Fluoranthene           380 U         UG/KG         2-Chloronaphthalene         380 U         UG/KG         Benzo(b)Fluoranthene           380 U         UG/						Phenanthrene
380 U         UG/KG         Naphthalene         380 U         UG/KG         Di-n-Butylphthalate           380 U         UG/KG         4-Chloroaniline         380 U         UG/KG         Fluoranthene           380 U         UG/KG         Hexachlorobutadiene         380 U         UG/KG         Pyrene           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Chrysene           380 U         UG/KG         2-4,5-Trichlorophenol         380 U         UG/KG         Di-n-Octylphthalate           380 U         UG/KG         2-4,5-Trichlorophenol         380 U         UG/KG         Benzo(b)Fluoranthene           380 U         UG/KG         2-Chloronaphthalene         380 U         UG/KG         Benzo(b)Fluoranthene           380 U         UG/KG		UG/KG		380 Ú	UG/KG	Anthracene
380 U         UG/KG         4-Chloroaniline         380 U         UG/KG         Fluoranthene           380 U         UG/KG         Hexachlorobutadiene         380 U         UG/KG         Pyrene           380 U         UG/KG         Caprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         4-Chloro-3-Methylphenol         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2-4,6-Trichlorophenol         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2,4,6-Trichlorophenol         380 U         UG/KG         Di-n-Octylphthalate           380 U         UG/KG         1,1-Biphenyl         380 U         UG/KG         Benzo(b)Fluoranthene           380 U         UG/KG         2-Chloronaphthalene         380 U         UG/KG         Benzo-a-Pyrene           380 U         UG/KG         2-Nitroaniline         380 U         UG/KG         Benzo-a-Pyrene           380 U         UG/KG         2,6-Dinitrotoluene         380 U         UG/KG         Dibenzo(a,h)Anthracene           380 U         UG/KG <td></td> <td></td> <td></td> <td></td> <td>UG/KG</td> <td></td>					UG/KG	
380 U         UG/KG         Hexachlorobutadiene         380 U         UG/KG         Pyrene           380 U         UG/KG         Acprolactam         380 U         UG/KG         Benzyl Butyl Phthalate           380 U         UG/KG         4-Chloro-3-Methylphenol         380 U         UG/KG         3,3'-Dichlorobenzidine           380 U         UG/KG         2-Methylnaphthalene         380 U         UG/KG         Benzo(a)Anthracene           380 U         UG/KG         2,4,6-Trichlorophenol         380 U         UG/KG         Chrysene           380 U         UG/KG         2,4,5-Trichlorophenol         380 U         UG/KG         Di-n-Octylphthalate           380 U         UG/KG         1,1-Biphenyl         380 U         UG/KG         Benzo(b)Fluoranthene           380 U         UG/KG         2-Chloronaphthalene         380 U         UG/KG         Benzo-a-Pyrene           380 U         UG/KG         2-Nitroaniline         380 U         UG/KG         Benzo-a-Pyrene           380 U         UG/KG         2,6-Dinitrotoluene         380 U         UG/KG         Dibenzo(a,h)Anthracene           380 U         UG/KG         2,6-Dinitrotoluene         380 U         UG/KG         Benzo(ghi)Perylene           380 U         UG/KG					UG/KG	Di-n-Butylphthalate
380 U UG/KG Caprolactam 380 U UG/KG 4-Chloro-3-Methylphenol 380 U UG/KG 4-Chloro-3-Methylphenol 380 U UG/KG 2-Methylnaphthalene 380 U UG/KG Benzo(a)Anthracene 380 U UG/KG Benzo(a)Anthracene 380 U UG/KG Chrysene 380 U UG/KG Chrysene 380 U UG/KG 2,4,6-Trichlorophenol 380 U UG/KG 2,4,5-Trichlorophenol 380 U UG/KG 2,4,5-Trichlorophenol 380 U UG/KG Din-Octylphthalate 380 U UG/KG 1,1-Biphenyl 380 U UG/KG Benzo(b)Fluoranthene 380 U UG/KG 2-Chloronaphthalene 380 U UG/KG 2-Nitroaniline 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dibenzo(a,h)Anthracene					UG/KG	Fluoranthene
380 U UG/KG 2-Methylnaphthalene 380 U UG/KG 380 U UG/KG Benzo(a)Anthracene 380 U UG/KG Benzo(a)Anthracene 380 U UG/KG Benzo(a)Anthracene 380 U UG/KG Benzo(a)Anthracene 380 U UG/KG Benzo(a)Anthracene 380 U UG/KG Chrysene 380 U UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG			i i i i i i i i i i i i i i i i i i i			
380 U       UG/KG       2-Methylnaphthalene       380 U       UG/KG       Benzo(a)Anthracene         380 U       UG/KG       Hexachlorocyclopentadiene (HCCP)       380 U       UG/KG       Chrysene         380 U       UG/KG       2,4,6-Trichlorophenol       380 U       UG/KG       Din-Octylphthalate         970 U       UG/KG       1,1-Biphenyl       380 U       UG/KG       Benzo(b)Fluoranthene         380 U       UG/KG       2-Chloronaphthalene       380 U       UG/KG       Benzo(b)Fluoranthene         970 U       UG/KG       2-Nitroaniline       380 U       UG/KG       Benzo-a-Pyrene         380 U       UG/KG       Dimethyl Phthalate       380 U       UG/KG       Benzo-a-Pyrene         380 U       UG/KG       2,6-Dinitrotoluene       380 U       UG/KG       Dibenzo(a,h)Anthracene         380 U       UG/KG       Acenaphthylene       380 U       UG/KG       Benzo(ghi)Perylene         380 U       UG/KG       3-Nitroaniline       14       % Moisture         380 U       UG/KG       2,4-Dinitrophenol						Benzyl Butyl Phthalate
380 UJ UG/KG Hexachlorocyclopentadiene (HCCP) 380 U UG/KG 2,4,6-Trichlorophenol 380 U UG/KG 2,4,5-Trichlorophenol 380 U UG/KG 2,4,5-Trichlorophenol 380 U UG/KG 1,1-Biphenyl 380 U UG/KG 2-Chloronaphthalene 380 U UG/KG 2-Chloronaphthalene 380 U UG/KG 2-Nitroaniline 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Benzo(b)Fluoranthene 380 U UG/KG Benzo(k)Fluoranthene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Benzo(a,h)Anthracene 380 U UG/KG Acenaphthylene 380 U UG/KG Acenaphthylene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene						
380 U UG/KG 2,4,5-Trichlorophenol 380 U UG/KG Di-n-Octylphthalate 380 U UG/KG Di-n-Octylphthalate 380 U UG/KG Di-n-Octylphthalate 380 U UG/KG Benzo(b)Fluoranthene 380 U UG/KG Benzo(b)Fluoranthene 380 U UG/KG Benzo(k)Fluoranthene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Indeno (1,2,3-cd) Pyrene 380 U UG/KG Acenaphthylene 380 U UG/KG Benzo(a,h)Anthracene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene						Benzo(a)Anthracene
970 U UG/KG 2,4,5-Trichlorophenol 380 U UG/KG Di-n-Octylphthalate 380 U UG/KG 1,1-Biphenyl 380 U UG/KG Benzo(b)Fluoranthene 380 U UG/KG 2-Chloronaphthalene 380 U UG/KG Benzo(k)Fluoranthene 970 U UG/KG 2-Nitroaniline 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 UJ UG/KG Indeno (1,2,3-cd) Pyrene 380 U UG/KG 2,6-Dinitrotoluene 380 UJ UG/KG Dibenzo(a,h)Anthracene 380 U UG/KG Acenaphthylene 380 UJ UG/KG Benzo(ghi)Perylene 970 U UG/KG 3-Nitroaniline 14 % Moisture  970 UR UG/KG 2,4-Dinitrophenol						
380 U UG/KG 1,1-Biphenyl 380 U UG/KG Benzo(b)Fluoranthene 380 U UG/KG 2-Chloronaphthalene 970 U UG/KG 2-Nitroaniline 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 UJ UG/KG Indeno (1,2,3-cd) Pyrene 380 U UG/KG 2,6-Dinitrotoluene 380 UJ UG/KG Dibenzo(a,h)Anthracene 380 U UG/KG Acenaphthylene 380 UJ UG/KG Benzo-(ghi)Perylene 970 U UG/KG 3-Nitroaniline 14 % Moisture 970 UR UG/KG 2,4-Dinitrophenol						bis(2-Ethylhexyl) Phthalate
380 U UG/KG 2-Chloronaphthalene 970 U UG/KG 2-Nitroaniline 380 U UG/KG Benzo(k)Fluoranthene 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 U UG/KG Indeno (1,2,3-cd) Pyrene 380 U UG/KG 2,6-Dinitrotoluene 380 U UG/KG Acenaphthylene 380 U UG/KG Acenaphthylene 380 U UG/KG Benzo(a,h)Anthracene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Benzo(ghi)Perylene						
970 U UG/KG 2-Nitroaniline 380 U UG/KG Benzo-a-Pyrene 380 U UG/KG Dimethyl Phthalate 380 UJ UG/KG Indeno (1,2,3-cd) Pyrene 380 U UG/KG 2,6-Dinitrotoluene 380 UJ UG/KG Dibenzo(a,h)Anthracene 380 U UG/KG Acenaphthylene 380 U UG/KG Benzo(ghi)Perylene 970 U UG/KG 3-Nitroaniline 14 % Moisture 970 UR UG/KG 2,4-Dinitrophenol						
380 U UG/KG Dimethyl Phthalate 380 U UG/KG 2,6-Dinitrotoluene 380 U UG/KG 2,6-Dinitrotoluene 380 U UG/KG Acenaphthylene 380 U UG/KG Acenaphthylene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 380 U UG/KG 2,4-Dinitrophenol						
380 U UG/KG 2,6-Dinitrotoluene 380 U UG/KG Acenaphthylene 380 U UG/KG Acenaphthylene 380 U UG/KG Benzo(ghi)Perylene 380 U UG/KG 3-Nitroaniline 380 U UG/KG Acenaphthene 380 U UG/KG Acenaphthene 970 UR UG/KG 2,4-Dinitrophenol						
380 U UG/KG Acenaphthylene 380 U UG/KG Benzo(ghi)Perylene 970 U UG/KG 3-Nitroaniline 14 % % Moisture 380 U UG/KG Acenaphthene 970 UR UG/KG 2,4-Dinitrophenol						Indeno (1,2,3-cd) Pyrene
970 U UG/KG 3-Nitroaniline 14 % % Moisture 380 U UG/KG Acenaphthene 970 UR UG/KG 2,4-Dinitrophenol			, · · · · · · · · · · · · · · · · ·			Dibenzo(a,h)Anthracene
380 U UG/KG Acenaphthene 970 UR UG/KG 2,4-Dinitrophenol				380 U	UG/KG	Benzo(ghi)Perylene
970 UR UG/KG 2,4-Dinitrophenol			· · · · · · · · · · · · · · · · · · ·	14	%	% Moisture
						•
970 UJ UG/KG 4-Nitrophenol						•
	970 UJ	UG/KG	4-Nitrophenol			
	,					

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 09:00

Requestor:

Ending:

Sample

6772 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS01SS /

Case No: 31635 MD No: 1XT2

Inorg Contractor: SENTIN

D No: 1XT2 Media: SURFACE SOIL (0" - 12")

Org Contractor: LIBRTY

**RESULTS UNITS** 1400 J\* UG/KG

**ANALYTE** 

12 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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Sample FY 2003 Project: 03-0474

**Extractables Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS13SS /

Flawood, MS

Case No: 31635

Inorg Contractor: SENTIN MD No: 1XT3

Org Contractor: LIBRTY D No: 1XT3 Media: SURFACE SOIL (0" - 12")

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE
- 1100 U	UG/KG	Benzaldehyde			1100 U	· UG/KG	Dibenzofuran
1100 U	UG/KG	Phenol			1100 U	UG/KG	2,4-Dinitrotoluene
1100 U	UG/KG	bis(2-Chloroethyl) Ether			1100 U	UG/KG	Diethyl Phthalate
1100 U	UG/KG	2-Chlorophenol			1100 U	UG/KG	Fluorene
1100 U	UG/KG	2-Methylphenol			1100 U	UG/KG	4-Chlorophenyl Phenyl Ether
1100 U	UG/KG	bis(2-Chloroisopropyl) Eth	er	•	2900 U	UG/KG	4-Nitroaniline
1100 U	UG/KG	Acetophenone		•	2900 U	UG/KG	2-Methyl-4,6-Dinitrophenol
1100 U	UG/KG	(3-and/or 4-)Methylphenol			1100 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
1100 U	UG/KG	n-Nitrosodi-n-Propylamine			NA	UG/KG	1,2,4,5-Tetrachlorobenzene
1100 U	UG/KG	Hexachloroethane		•	1100 U	UG/KG	4-Bromophenyl Phenyl Ether
1100 U	UG/KG	Nitrobenzene		•	1100 U	UG/KG	Hexachlorobenzene (HCB)
1100 U	UG/KG	Isophorone			1100 U	UG/KG	Atrazine
1100 U	UG/KG	2-Nitrophenol			2900 U	- UG/KG	Pentachlorophenol
1100 U	UG/KG	2,4-Dimethylphenol			1600	UG/KG	Phenanthrene
1100 U	UG/KG	bis(2-Chloroethoxy)Methar	ne	. '	1600	UG/KG	Anthracene
1100 U	UG/KG	2,4-Dichlorophenol	•		540 J	UG/KG	Carbazole
1100 U	UG/KG	Naphthalene	•		1100 U	UG/KG	Di-n-Butylphthalate
1100 U	UG/KG	4-Chloroaniline			11000	UG/KG	Fluoranthene
1100 U -	UG/KG	Hexachlorobutadiene			6800	UG/KG	Pyrene
1100 U	UG/KG	Caprolactam	•		1100 UJ	UG/KG	Benzyl Butyl Phthalate
1100 U	UG/KG	4-Chloro-3-Methylphenol			1100 UJ	UG/KG	3,3'-Dichlorobenzidine
1100 U	UG/KG	2-Methylnaphthalene			4300	· UG/KG	Benzo(a)Anthracene
1100 U	UG/KG	Hexachlorocyclopentadien	ie (HCCP)		6200	UG/KG	Chrysene
1100 U	UG/KG	2,4,6-Trichlorophenol		•	1100 U	UG/KG	bis(2-Ethylhexyl) Phthalate
2900 U	UG/KG	2,4,5-Trichlorophenol			1100 U	UG/KG	Di-n-Octylphthalate
1100 U	UG/KG	1,1-Biphenyl		•	7300 J	UG/KG	Benzo(b)Fluoranthene
1100 U	UG/KG	2-Chloronaphthalene			5400 J	UG/KG	Benzo(k)Fluoranthene
2900 U	UG/KG	2-Nitroaniline			3100	UG/KG	Benzo-a-Pyrene
1100 U	UG/KG	Dimethyl Phthalate			3100	UG/KG	Indeno (1,2,3-cd) Pyrene
1100 U	UG/KG	2,6-Dinitrotoluene			1100	UG/KG	Dibenzo(a,h)Anthracene
1000 J	UG/KG	Acenaphthylene			2000	UG/KG	Benzo(ghi)Perylene
2900 U	UG/KG	3-Nitroaniline	•		13	%	% Moisture
1100 U	UG/KG	Acenaphthene					
2900 U	UG/KG	2,4-Dinitrophenol					
2900 U	UG/KG	4-Nitrophenol					
				•			•

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474 Sample 6773 FY 2003

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS13SS /

MD No: 1XT3

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XT3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:15

Ending:

UNITS	ANALYTE
UG/KG	PHENANTHRENE, 1-METHYL-
UG/KG	9,10-ANTHRACENEDIONE
UG/KG	PHENANTHRENE, 2,3-DIMETHYL-
UG/KG	CYCLOPENTA (DEF) PHENANTHRENONE
UG/KG	BENZO [B] NAPHTHO [2,3-D] FURAN
UG/KG	11H-BENZO [B] FLUORENE
UG/KG	PYRENE, 1-METHYL-
UG/KG	UNKNOWN PAH
UG/KG	PYRENE, 2-METHYL-
UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
UG/KG	CHRYSENE, 5-METHYL-
UG/KG	5,12-NAPHTAHCENEDIONE
UG/KG	2,2'-BINAPHTHALANE
UG/KG	6 UNKNOWN PAHS
UG/KG	12 UNKNOWN COMPOUNDS
	UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Media: SUBSURFACE SOIL (> 12")

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise Project: 03-0474 FY 2003 Sample Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 13:25 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS13SB / MD No: 1XT4

Org Contractor: LIBRTY

	<del></del>				
RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
390 U	UG/KG	Benzaldehyde	390 U	UG/KG	Dibenzofuran
390 Ư	UG/KG	Phenol	390 U	UG/KG	2,4-Dinitrotoluene
390 U	UG/KG	bis(2-Chloroethyl) Ether	390 U	UG/KG	Diethyl Phthalate
390 U	UG/KG	2-Chlorophenol	390 U	UG/KG	Fluorene
390 U	UG/KG	2-Methylphenol	390 U	UG/KG	4-Chlorophenyl Phenyl Ether
390 U	UG/KG	bis(2-Chloroisopropyl) Ether	980 U	UG/KG	4-Nitroaniline
390 U	UG/KG	Acetophenone	980 U	UG/KG	2-Methyl-4,6-Dinitrophenol
390 U	UG/KG	(3-and/or 4-)Methylphenol	390 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
390 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
390 U	UG/KG	Hexachloroethane	390 U	UG/KG	4-Bromophenyl Phenyl Ether
390 U	UG/KG	Nitrobenzene	390 U	UG/KG	Hexachlorobenzene (HCB)
. 390 U	UG/KG	Isophorone	390 U	UG/KG	Atrazine
390 U	UG/KG	2-Nitrophenol	980 U	UG/KG	Pentachlorophenol
390 U	UG/KG	2,4-Dimethylphenol	390 U	UG/KG	Phenanthrene
390 U	'UG/KG	bis(2-Chloroethoxy)Methane	390 U	UG/KG	Anthracene
390 U	UG/KG	2,4-Dichlorophenol	390 U	UG/KG	Carbazole
390 U	UG/KG	Naphthalene	390 U	UG/KG	Di-n-Butylphthalate
390 U	UG/KG	4-Chloroaniline	130 J	UG/KG	Fluoranthene
390 U	UG/KG	Hexachlorobutadiene	89 J	UG/KG	Pyrene
390 U	UG/KG	Caprolactam	390 UJ	UG/KG	Benzyl Butyl Phthalate
390 U	UG/KG	4-Chloro-3-Methylphenol	390 UJ	UG/KG	3,3'-Dichlorobenzidine
390 U	UG/KG	2-Methylnaphthalene	56 J	UG/KG	Benzo(a)Anthracene
390 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	91 J	UG/KG	Chrysene
390 U	UG/KG	2,4,6-Trichlorophenol	390 U	UG/KG	bis(2-Ethylhexyl) Phthalate
980 U	UG/KG	2,4,5-Trichlorophenol	390 U	UG/KG	Di-n-Octylphthalate
390 U	UG/KG	1,1-Biphenyl	160 J	UG/KG	Benzo(b)Fluoranthene
390 U	UG/KG	2-Chloronaphthalene	170 J	UG/KG	Benzo(k)Fluoranthene
980 U	UG/KG	2-Nitroaniline	390 U	UG/KG	Benzo-a-Pyrene
390 U -	UG/KG	Dimethyl Phthalate	51 J	UG/KG	Indeno (1,2,3-cd) Pyrene
390 U	UG/KG	2,6-Dinitrotoluene	390 U	UG/KG	Dibenzo(a,h)Anthracene
390 Ù	UG/KG	Acenaphthylene	390 U	UG/KG	Benzo(ghi)Perylene
980 U	UG/KG	3-Nitroaniline	15	%	% Moisture
390 U	UG/KG	Acenaphthene			
980 U	UG/KG	2,4-Dinitrophenol			•
980 U	UG/KG	4-Nitrophenol			
•					

D No: 1XT4

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6774 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 13:25

Ending:

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF Id/Station: GS13SB / Case No: 31635

MD No: 1XT4

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

UG/KG

D No: 1XT4

Org Contractor: LIBRTY

**RESULTS UNITS** 

1000 J

**ANALYTE** 

2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

Produced by: Goddard, Denise Project: 03-0474 Sample 6775 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 11:35 Program: SF Case No: 31635 Ending: ld/Station: GS04SD / Inorg Contractor: SENTIN MD No: 1XT5 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XT5 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
130 J	UG/KG	Benzaldehyde	900 U	UG/KG	Dibenzofuran
900 U	UG/KG	Phenol	900 U	UG/KG	2,4-Dinitrotoluene
900 U	UG/KG	bis(2-Chloroethyl) Ether	900 U	UG/KG	Diethyl Phthalate
900 U	UG/KG	2-Chlorophenol	900 U	UG/KG	Fluorene
900 U	UG/KG	2-Methylphenol	900 U	UG/KG	4-Chlorophenyl Phenyl Ether
900 U	UG/KG	bis(2-Chloroisopropyl) Ether	2300 U	UG/KG	4-Nitroaniline
900 U	UG/KG	Acetophenone	2300 U	UG/KG	2-Methyl-4,6-Dinitrophenol
900 U	UG/KG	(3-and/or 4-)Methylphenol	900 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
900 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
900 U	UG/KG	Hexachloroethane	900 U	UG/KG	4-Bromophenyl Phenyl Ether
900 U -	UG/KG	Nitrobenzene	900 U	UG/KG	Hexachlorobenzene (HCB)
900 U	UG/KG	Isophorone	900 U	UG/KG	Atrazine
900 U	UG/KG	2-Nitrophenol	2300 U	UG/KG	Pentachlorophenol
900 U	UG/KG	2,4-Dimethylphenol	900 U	UG/KG	Phenanthrene
900 U	UG/KG	bis(2-Chloroethoxy)Methane	900 U	UG/KG	Anthracene
900 U	UG/KG	2,4-Dichlorophenol	900 U	UG/KG	Carbazole
3.0 J	UG/KG	Naphthalene	900 U	UG/KG	Di-n-Butylphthalate
900 U	UG/KG	4-Chloroaniline	900 U	UG/KG	Fluoranthene
900 U	UĠ/KG	Hexachlorobutadiene	120 J	UG/KG	Pyrene
900 U	UG/KG	Caprolactam	900 UJ	UG/KG	Benzyl Butyl Phthalate
900 U	UG/KG	4-Chloro-3-Methylphenol	900 UJ	UG/KG	3,3'-Dichlorobenzidine
900 U	UG/KG	2-Methylnaphthalene	900 U	UG/KG	Benzo(a)Anthracene
900 U	UG/KĠ	Hexachlorocyclopentadiene (HCCP)	900 U	UG/KG	Chrysene
900 U	UG/KG	2,4,6-Trichlorophenol	900 U	UG/KG	bis(2-Ethylhexyl) Phthalate
2300 U	UG/KG	2,4,5-Trichlorophenol	900 U	UG/KG	Di-n-Octylphthalate
900 U	UG/KG	1,1-Biphenyl	900 U	UG/KG	Benzo(b)Fluoranthene
900 U	UG/KG	2-Chloronaphthalene	900 U	UG/KG	Benzo(k)Fluoranthene
2300 U	UG/KG	2-Nitroaniline	900 U	UG/KG	Benzo-a-Pyrene
900 U	UG/KG	Dimethyl Phthalate	900 U	UG/KG	Indeno (1,2,3-cd) Pyrene
900 U	UG/KG	2,6-Dinitrotoluene	900 U	UG/KG	Dibenzo(a,h)Anthracene
900 U	UG/KG	Acenaphthylene	900 U	UG/KG	Benzo(ghi)Perylene
2300 U	UG/KG	3-Nitroaniline	78	%	% Moisture
900 U	UG/KG	Acenaphthene		•	
2300 U	UG/KG	2,4-Dinitrophenol			
2300 U	UG/KG	4-Nitrophenol			
•					

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

EPA - REGION IV SESD, ATHENS, GA

Production Date: 06/20/2003 14:10

Sample

Project: 03-0474 6775 FY 2003

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS04SD / Media: SEDIMENT

Case No: 31635

MD No: 1XT5

D No: 1XT5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:35

Ending:

RESULTS UNITS **ANALYTE** 

UG/KG

1600 NJ UG/KG **CHOLESTANOL** 

1700 NJ UG/KG 9200 J UG/KG ERGOST-5-EN-3-OL, (3.BETA.)-

1800 J UG/KG UG/KG 2200 NJ

54000 J

UNKNOWN ALCOHOL UNKNOWN KETONE STIGMAST-4-EN-3-ONE

24 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Facility: Gulf States Creosoting

6775 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 11:35

Ending:

Program: SF Id/Station: GS04SD /

**SPECIFIED TESTS** 

Flowood, MS

Case No: 31635

Inorg Contractor: SENTIN

Media: SEDIMENT

MD No: 1XT5 D No: 1XT5

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2 J	UG/KG	2-Methylnaphthalene
3 J	UG/KG	Naphthalene
9 .	UG/KG	Acenaphthylene
5	UG/KG	Acenaphthene
6	UG/KG	Fluorene
11	UG/KG	Phenanthrene
34	UG/KG	Anthracene
50	UG/KG	Fluoranthene
33	UG/KG	Pyrene
19	UG/KG	Benzo(a)Anthracene
36	UG/KG	Chrysene
NA	UG/KG	Benzo(b)Fluoranthene
NA	UG/KG	Benzo(k)Fluoranthene
1,7	.UG/KG	Benzo-a-Pyrene
NA	UG/KG	Indeno (1,2,3-cd) Pyrene
6	UG/KG	Dibenzo(a,h)Anthracene
NA	UG/KG	Benzo(ghi)Perylene
49 U	UG/KG	Pentachlorophenol
•		-

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise 6776 FY 2003 Project: 03-0474 Sample Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 12:15 Program: SF Case No: 31635 Endina: Id/Station: GS05SD / MD No: 1XT6 Inorg Contractor: SENTIN D No: 1XT6 Org Contractor: LIBRTY Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS RESULTS UNITS **ANALYTE RESULTS UNITS ANALYTE** 430 U UG/KG Benzaldehyde 430 U UG/KG Dibenzofuran 430 U UG/KG Phenol 430 U UG/KG 2.4-Dinitrotoluene 430 U UG/KG bis(2-Chloroethyl) Ether 430 U UG/KG Diethyl Phthalate 430 U UG/KG 2-Chlorophenol 430 U UG/KG Fluorene 430 U UG/KG 2-Methylphenol 430 U UG/KG 4-Chlorophenyl Phenyl Ether 430 U bis(2-Chloroisopropyl) Ether 1100 UJ UG/KG UG/KG 4-Nitroaniline 430 U UG/KG Acetophenone 1100 U UG/KG 2-Methyl-4,6-Dinitrophenol 430 U UG/KG (3-and/or 4-)Methylphenol 430 U UG/KG n-Nitrosodiphenylamine/Diphenylamine 430 U UG/KG n-Nitrosodi-n-Propylamine NA UG/KG 1,2,4,5-Tetrachlorobenzene 430 U UG/KG Hexachloroethane 430 U UG/KG 4-Bromophenyl Phenyl Ether 430 U Nitrobenzene UG/KG 430 UJ UG/KG Hexachlorobenzene (HCB) 430 U UG/KG Isophorone 430 U UG/KG Atrazine 430 U UG/KG 2-Nitrophenol 1100 UJ UG/KG Pentachlorophenol 430 U UG/KG 2.4-Dimethylphenol 430 U UG/KG Phenanthrene 430 U UG/KG bis(2-Chloroethoxy)Methane 430 U UG/KG Anthracene 430 U UG/KG 2,4-Dichlorophenol 430 U UG/KG Carbazole Naphthalene 430 U UG/KG 430 U UG/KG Di-n-Butylphthalate 430 U UG/KG 4-Chloroaniline 430 U UG/KG Fluoranthene 430 U UG/KG Hexachlorobutadiene 430 U UG/KG Pyrene 430 U UG/KG Caprolactam 430 U UG/KG Benzyl Butyl Phthalate 430 U UG/KG 4-Chloro-3-Methylphenol 430 U UG/KG 3,3'-Dichlorobenzidine 430 U UG/KG 2-Methylnaphthalene 430 U UG/KG Benzo(a)Anthracene 430 UJ UG/KG Hexachlorocyclopentadiene (HCCP) 430 U UG/KG Chrysene 430 U UG/KG 2,4,6-Trichlorophenol 430 U UG/KG bis(2-Ethylhexyl) Phthalate 1100 U UG/KG 2,4,5-Trichlorophenol 430 U UG/KG Di-n-Octylphthalate 430 U UG/KG 1.1-Biphenvl 430 U UG/KG Benzo(b)Fluoranthene 430 U UG/KG 2-Chloronaphthalene 430 U UG/KG Benzo(k)Fluoranthene 1100 U UG/KG 2-Nitroaniline 430 U UG/KG Benzo-a-Pyrene 430 U UG/KG Dimethyl Phthalate 430 UJ UG/KG Indeno (1,2,3-cd) Pyrene 430 U UG/KG 2,6-Dinitrotoluene 430 UJ UG/KG Dibenzo(a,h)Anthracene 430 U UG/KG Acenaphthylene 430 U UG/KG Benzo(ghi)Perylene 1100 U UG/KG 3-Nitroaniline 54 % Moisture 430 U UG/KG Acenaphthene 2,4-Dinitrophenol 1100 UR UG/KG 1100 UJ UG/KG 4-Nitrophenol

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyzed in replicate. Reported value is "average" of replicates.

## **EPA - REGION IV SESD, ATHENS, GA**

EXTRACTABLES SAMPLE ANALYSIS

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6776 FY **2003** 

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS05SD / Media: SEDIMENT

Case No: 31635

MD No: 1XT6

D No: 1XT6

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Ending:

Requestor:

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 12:15

**RESULTS UNITS ANALYTE** 

1200 NJ UG/KG 500 J UG/KG 2500 NJ UG/KG

BENZENE, 1-METHYL-2-ISOPROPYL **UNKNOWN CARBOXYLIC ACID** PHENANTHRENONE DERIVATIVE STIGMAST-4-EN-3-ONE

800 NJ UG/KG 25 UNKNOWN COMPOUNDS 140000 J UG/KG

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6776 FY 2003

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS05SD / Media: SEDIMENT

Flowood, MS

Case No: 31635 MD No: 1XT6

D No: 1XT6

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.8 J	UG/KG	2-Methylnaphthalene
0.9 J	UG/KG	Naphthalene
2 U	UG/KG	Acenaphthylene
0.9 J	UG/KG	Acenaphthene
1 J	UG/KG	Fluorene
23 U	UG/KG	Phenanthrene
<b>7</b> ·	UG/KG	Anthracene
34	UG/KG	Fluoranthene
27	UG/KG	Pyrene
12	UG/KG	Benzo(a)Anthracene
26	UG/KG	Chrysene
NA	UG/KG	Benzo(b)Fluoranthene
NA	UG/KG	Benzo(k)Fluoranthene
16	UG/KG	Benzo-a-Pyrene
NA	UG/KG	Indeno (1,2,3-cd) Pyrene
4	UG/KG	Dibenzo(a,h)Anthracene
NA	UG/KG	Benzo(ghi)Perylene
- 23 U	UG/KG	Pentachlorophenol
		·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Sample FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 12:45 Program: SF Case No: 31635 Ending: - Inorg Contractor: SENTIN Id/Station: GS07SD / MD No: 1XT7 Org Contractor: LIBRTY D No: 1XT7 Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
			330 U		
330 U	UG/KG	Benzaldehyde		UG/KG	Dibenzofuran
330 U	UG/KG	Phenol	330 U	UG/KG	2,4-Dinitrotoluene
330 U	UG/KG	bis(2-Chloroethyl) Ether	330 U	UG/KG	Diethyl Phthalate
330 U	UG/KG	2-Chlorophenol	330 U	UG/KG	Fluorene
330 U	UG/KG	2-Methylphenol	330 U	UG/KG	4-Chlorophenyl Phenyl Ether
330 U	UG/KG	bis(2-Chloroisopropyl) Ether	830 UJ	UG/KG	4-Nitroaniline
330 U	UG/KG	Acetophenone	830 U	UG/KG	2-Methyl-4,6-Dinitrophenol
. 330 U	UG/KG	(3-and/or 4-)Methylphenol	330 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
330 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
330 U	UG/KG	Hexachloroethane	330 U	UG/KG	4-Bromophenyl Phenyl Ether
330 U	UG/KG	Nitrobenzene	330 UJ	UG/KG	Hexachlorobenzene (HCB)
330 U -	UG/KG	Isophorone	330 U	UG/KG	Atrazine
330 U	UG/KG	2-Nitrophenol	830 UJ	UG/KG	Pentachlorophenol
330 U	UG/KG	2,4-Dimethylphenol	330 U	UG/KG	Phenanthrene
330 U	UG/KG	bis(2-Chloroethoxy)Methane	330 U	UG/KG	Anthracene
330 U	UG/KG	2,4-Dichlorophenol	330 U	UG/KG	Carbazole
330 U	UG/KG	Naphthalene	330 U	UG/KG	Di-n-Butylphthalate
330 U	UG/KG	4-Chloroaniline	330 U	UG/KG	Fluoranthene
330 U	UG/KG	Hexachlorobutadiene	330 U	UG/KG	Pyrene
330 U	UG/KG	Caprolactam	330 U	UG/KG	Benzyl Butyl Phthalate
330 U	UG/KG	4-Chloro-3-Methylphenal	330 U	UG/KG	3,3'-Dichlorobenzidine
330 U	UG/KG	2-Methylnaphthalene	330 U	UG/KG	Benzo(a)Anthracene
330 UJ	UG/KG	Hexachlorocyclopentadiene (HCCP)	330 U ·	UG/KG	Chrysene
330 U	UG/KG	2,4,6-Trichlorophenol	330 U	UG/KG	bis(2-Ethylhexyl) Phthalate
830 U	UG/KG	2,4,5-Trichlorophenol	330 U	UG/KG	Di-n-Octylphthalate
330 U	UG/KG	1,1-Biphenyl	330 U	UG/KG	Benzo(b)Fluoranthene
330 U	UG/KG	2-Chloronaphthalene	330 U	UG/KG	Benzo(k)Fluoranthene
830 U	UG/KG	2-Nitroaniline	330 U	UG/KG	Benzo-a-Pyrene
330 U	UG/KG	Dimethyl Phthalate	330 UJ	UG/KG	Indeno (1,2,3-cd) Pyrene
330 U	UG/KG	2,6-Dinitrotoluene	330 UJ	UG/KG	Dibenzo(a,h)Anthracene
330 U	UG/KG	Acenaphthylene	330 U	UG/KG	Benzo(ghi)Perylene
830 U	UG/KG	3-Nitroaniline	18	%	% Moisture
	UG/KG	Acenaphthene	- <del>-</del>		· · · · · · · · · · · · · · · · · ·
830 UR	UG/KG	2,4-Dinitrophenol			
830 UJ	UG/KG	4-Nitrophenol			
			•		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 12:45

Requestor:

Ending:

Sample

6777

FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Id/Station: GS07SD /

Program: SF

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Flowood, MS

MD No: 1XT7 D No: 1XT7

**RESULTS UNITS** 

Media: SEDIMENT

69 J UG/KG **ANALYTE** 

UNKNOWN COMPOUND

Data Reported as Identified by CLP Lab - IDs Not Verified

# **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6777 FY 2003

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Media: SEDIMENT

Id/Station: GS07SD /

Case No: 31635

MD No: 1XT7 D No: 1XT7

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:45

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.4 J	UG/KG	2-Methylnaphthalene
0.4 J	UG/KG	Naphthalene
0.8 U	UG/KG	Acenaphthylene
0.9 U	UG/KG	Acenaphthene
0.8 U	UG/KG	Fluorene
0.9 U	UG/KG	Phenanthrene
0.8 U	UG/KG	Anthracene
2 U	UG/KG	Fluoranthene
2 U	UG/KG	Pyrene
0.8 U	UG/KG	Benzo(a)Anthracene
1 U	UG/KG	Chrysene
NA	UG/KG	Benzo(b)Fluoranthene
NA	UG/KG	Benzo(k)Fluoranthene
0.8 U	UG/KG	Benzo-a-Pyrene
NA	UG/KG	Indeno (1,2,3-cd) Pyrene
0.8 U	UG/KG	Dibenzo(a,h)Anthracene
NA	UG/KG	Benzo(ghi)Perylene
13 U	UG/KG	Pentachlorophenol
		į

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise 6778 FY 2003 Project: 03-0474 Sample Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:55 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS23SS / MD No: 1XT8 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT8 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE		RESULTS	UNITS	ANALYTE
450 U	UG/KG	Benzaldehyde		450 U	UG/KG	Dibenzofuran
450 U	UG/KG	Phenol		450 U	UG/KG	2,4-Dinitrotoluene
450 U	UG/KG	bis(2-Chloroethyl) Ether		450 U	UG/KG	Diethyl Phthalate
450 U	UG/KG	2-Chlorophenol		450 U	UG/KG	Fluorene
450 U	UG/KG	2-Methylphenol	•	450 U	UG/KG	4-Chlorophenyl Phenyl Ether
450 U	UG/KG	bis(2-Chloroisopropyl) Ether		1100 UJ	UG/KG	4-Nitroaniline
450 U	UG/KG	Acetophenone		1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
450 U	UG/KG	(3-and/or 4-)Methylphenol		450 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
450 U	UG/KG	n-Nitrosodi-n-Propylamine		· NA	UG/KG	1,2,4,5-Tetrachlorobenzene
450 U	UG/KG	Hexachloroethane		450 U	UG/KG	4-Bromophenyl Phenyl Ether
450 U	UG/KG	Nitrobenzene	•	450 UJ	UG/KG	Hexachlorobenzene (HCB)
450 U	UG/KG	Isophorone		` 450 U	UG/KG	Atrazine
450 Ư	UG/KG	2-Nitrophenol		1100 UJ	UG/KG	Pentachlorophenol
450 U	UG/KG	2,4-Dimethylphenol		59 J	UG/KG	Phenanthrene
450 U	UG/KG	bis(2-Chloroethoxy)Methane		430 U	UG/KG	Anthracene
450 U	UG/KG	2,4-Dichlorophenol	•	450 U	UG/KG	Carbazole
450 U	UG/KG	Naphthalene		450 U	UG/KG	Di-n-Butylphthalate
450 U	UG/KG	4-Chloroaniline		520	UG/KG	Fluoranthene
450 U	UG/KG	Hexachlorobutadiene		450	UG/KG	Pyrene
450 U	UG/KG	Caprolactam		450 U	UG/KG	Benzyl Butyl Phthalate
450 U	UG/KG	4-Chloro-3-Methylphenol		450 U	UG/KG	3,3'-Dichlorobenzidine
450 U	UG/KG	2-Methylnaphthalene		430 J	UG/KG	Benzo(a)Anthracene
450 UJ	UG/KG	Hexachlorocyclopentadiene (HCCP)		510	UG/KG	Chrysene
450 U	UG/KG	2,4,6-Trichlorophenol	•	450 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol		450 U	UG/KG	Di-n-Octylphthalate
450 U	UG/KG	1,1-Biphenyl			UG/KG	Benzo(b)Fluoranthene
450 U	UG/KG	2-Chloronaphthalene		520	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline		380 J	UG/KG	Benzo-a-Pyrene
450 U	UG/KG	Dimethyl Phthalate		340 J	UG/KG	Indeno (1,2,3-cd) Pyrene
450 U	UG/KG	2,6-Dinitrotoluene		120 J	UG/KG	Dibenzo(a,h)Anthracene
66 J	UG/KG	Acenaphthylene		180 J	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline		26	%	% Moisture
450 U	UG/KG	Acenaphthene				
1100 UR	UG/KG	2,4-Dinitrophenol				
1100 UJ	UG/KG	4-Nitrophenol				

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value,

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

Sample 6778 FY 2003 Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

140 NJ

0,1000

Id/Station: GS23SS /

Media: SURFACE SOIL (0" - 12")

UG/KG

Case No: 31635

MD No: 1XT8 D No: 1XT8 Inorg Contractor: SENTIN

Org Contractor: LIBRTY

J

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 13:55

Requestor:

Ending:

RESULTS UNITS ANALYTE

CYCLOPENTA (DEF) PHENANTHRENONE

130 NJ UG/KG 210 NJ UG/KG PYRENE, 2-METHYL-

210 NJ UG/KG 240 NJ UG/KG 7H-BENZ [DE] ANTHRACEN-7-ONE BENZ [A] ANTHRACENE. 7-METHYL-

210 NJ UG/KG

5. 12-NAPHTHACENEDIONE

240 NJ UG/KG

BENZO [E] PYRENE

440 NJ UG/KG 110 NJ UG/KG BENZO [J] FLUORANTHENE BENZ [E] ACEPHENANTHRYLENE

3600 J UG/KG 170 NJ UG/KG 21 UNKNOWN COMPOUNDS STIGMAST-4-EN-3-ONE

Data Reported as Identified by CLP Lab - IDs Not Verified

990 UJ

UG/KG

4-Nitrophenol

Produced by: Goddard, Denise 6779 FY 2003 Project: 03-0474 Sample Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 13:40 Program: SF Case No: 31635 Ending: Id/Station: GS15SS / MD No: 1XT9 Inorg Contractor: SENTIN Org Contractor: LIBRTY D No: 1XT9 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS ANALYTE** 390 U UG/KG Dibenzofuran 390 U UG/KG Benzaldehyde 390 U UG/KG Phenol 390 U UG/KG 2.4-Dinitrotoluene 390 U bis(2-Chloroethyl) Ether 390 U UG/KG Diethyl Phthalate UG/KG 390 U 390 U UG/KG Fluorene UG/KG 2-Chlorophenol 4-Chlorophenyl Phenyl Ether 390 U 2-Methylphenol 390 U UG/KG UG/KG 390 U bis(2-Chloroisopropyl) Ether 990 UJ UG/KG 4-Nitroaniline UG/KG

390 U UG/KG Acetophenone 390 U UG/KG (3-and/or 4-)Methylphenol 390 U UG/KG n-Nitrosodi-n-Propylamine 390 U UG/KG Hexachloroethane 390 U Nitrobenzene UG/KG 390 U UG/KG Isophorone 390 U ·UG/KG 2-Nitrophenol 390 U UG/KG 2.4-Dimethylphenol 390 U UG/KG bis(2-Chloroethoxy)Methane 390 U UG/KG 2,4-Dichlorophenol. 390 U UG/KG Naphthalene 390 U UG/KG 4-Chloroaniline 390 U UG/KG Hexachlorobutadiene 390 U UG/KG Caprolactam 390 U 4-Chloro-3-Methylphenol UG/KG 390 U UG/KG 2-Methylnaphthalene 390 UJ UG/KG Hexachlorocyclopentadiene (HCCP) 390 U UG/KG 2,4,6-Trichlorophenol 990 U UG/KG 2.4.5-Trichlorophenol 390 U UG/KG 1,1-Biphenyl 390 U UG/KG 2-Chloronaphthalene 990 U UG/KG 2-Nitroaniline 390 U Dimethyl Phthalate UG/KG 2.6-Dinitrotoluene 390 U UG/KG 250 J UG/KG Acenaphthylene 990 U UG/KG 3-Nitroaniline 390 U UG/KG Acenaphthene 990 UR UG/KG 2.4-Dinitrophenol

990 U UG/KG 2-Methyl-4,6-Dinitrophenol 390 U UG/KG n-Nitrosodiphenylamine/Diphenylamine NA UG/KG 1,2,4,5-Tetrachlorobenzene 390 U UG/KG 4-Bromophenyl Phenyl Ether UG/KG Hexachlorobenzene (HCB) 390 UJ 390 U -UG/KG Atrazine 990 UJ UG/KG Pentachlorophenol 1100 UG/KG Phenanthrene 370 J UG/KG Anthracene 250 J UG/KG Carbazole 390 U UG/KG Di-n-Butylphthalate 3100 UG/KG Fluoranthene 2200 UG/KG Pvrene 390 U UG/KG Benzyl Butyl Phthalate 390 U UG/KG 3,3'-Dichlorobenzidine 690 UG/KG Benzo(a)Anthracene 1700 UG/KG Chrysene 390 U UG/KG bis(2-Ethylhexyl) Phthalate 390 U UG/KG Di-n-Octylphthalate Benzo(b)Fluoranthene 1500 UG/KG 1100 UG/KG Benzo(k)Fluoranthene 650 UG/KG Benzo-a-Pyrene 740 J UG/KG Indeno (1,2,3-cd) Pyrene 200 J Dibenzo(a,h)Anthracene UG/KG 360 J UG/KG Benzo(ghi)Perylene 16 % % Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

EPA - REGION IV SESD. ATHENS. GA

Production Date: 06/20/2003 14:10

Sample

6779 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS15SS /

MD No: 1XT9

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

Org Contractor: LIBRTY D No: 1XT9

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:40

Ending:

**RESULTS UNITS ANALYTE** 

700 NJ UG/KG 350 NJ UG/KG 9,10-ANTHRACENEDIONE 11H-BENZO [B] FLUORENE

440 NJ UG/KG 7H-BENZ [DE] ANTHRACEN-7-ONE

300 NJ UG/KG

CHRYSENE, 1-METHYL-

BENZ (A) ANTHRACENE-7, 12-DIONE 260 NJ UG/KG **4 UNKNOWN PAHS** 

UG/KG 2000 J 12000 J UG/KG

19 UNKNOWN COMPOUNDS

1,2:3,4-DIBENZPYRENE 190 NJ UG/KG

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample 6780 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 13:50 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN ld/Station: GS15SB / MD No: 1XW0 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XW0 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
420 U	UG/KG	Benzaldehyde	420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	420 U	UG/KG	Diethyl Phthalate
420 U	UG/KG	2-Chlorophenol	420 U	UG/KG	Fluorene
420 U	UG/KG	2-Methylphenol	420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 UJ	UG/KG	4-Nitroaniline
420 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
420 U	UG/KG	(3-and/or 4-)Methylphenol	420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane	420 U	UG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	420 UJ	UG/KG	Hexachlorobenzene (HCB)
420 U	UG/KG	Isophorone	.420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol	1100 UJ	UG/KG	Pentachlorophenol
420 U	UG/KG	2,4-Dimethylphenol	420 U	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Methane	420 U	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol	420 U	UG/KG	Carbazole
420 U	UG/KG	Naphthalene	420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline	· 420 U	UG/KG	Fluoranthene
420 U	UG/KG	Hexachlorobutadiene	420 U	UG/KG	Pyrene
420 U	UG/KG	Caprolactam	420 U	UG/KG	Benzyl Butyl Phthalate
420 U	UG/KG	4-Chloro-3-Methylphenal	420 U	UG/KG	3,3'-Dichlorobenzidine
420 U	UG/KG	2-Methylnaphthalene	420 U	UG/KG	Benzo(a)Anthracene
420 UJ	UG/KG	Hexachlorocyclopentadiene (HCCP)	420 U	UG/KG	Chrysene
420 U	UG/KG	2,4,6-Trichlorophenol	420 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	420 U	UG/KG	Di-n-Octylphthalate
420 U	UG/KG	1,1-Biphenyl	420 U	UG/KG	Benzo(b)Fluoranthene
420 U	UG/KG	2-Chloronaphthalene	420 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	420 U	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate	420 UJ	UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2,6-Dinitrotoluene	420 UJ	UG/KG	Dibenzo(a,h)Anthracene
420 U	UG/KG	Acenaphthylene	420 U	UG/KG	Benzo(ghi)Perylene
1100 U -	UG/KG	3-Nitroaniline	21	%	% Moisture
420 U	UG/KG	Acenaphthene	•		. •
1100 UR	UG/KG	2,4-Dinitrophenol			
1100 UJ	UG/KG	4-Nitrophenol	Α.	•	
		1			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6780 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS15SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

Inorg Contractor: SENTIN MD No: 1XW0 D No: 1XW0

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:50

Ending:

RESULTS UNITS 310 J UG/KG

**ANALYTE** 

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample 6781 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:13 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS02SS / MD No: 1XW1

Media: SURFACE SOIL (0" - 12") D No: 1XW1 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
430 U	UG/KG	Benzaldehyde	430 U	UG/KG	Dibenzofuran
430 U	UG/KG	Phenol	430 U	UG/KG	2,4-Dinitrotoluene
430 U	UG/KG	bis(2-Chloroethyl) Ether	430 U	UG/KG	Diethyl Phthalate
430 U	UG/KG	2-Chlorophenol	430 U	UG/KG	Fluorene
430 U	UG/KG	2-Methylphenol	430 U	UG/KG	4-Chlorophenyl Phenyl Ether
430 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 UJ	UG/KG	4-Nitroaniline
: 430 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
430 U	UG/KG	(3-and/or 4-)Methylphenol	430 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
430 U	UG/KG	n-Nitrosodi-n-Propylamine	. NA	UG/KG	1,2,4,5-Tetrachlorobenzene
430 U	UG/KG	Hexachloroethane	430 U	UG/KG	4-Bromophenyl Phenyl Ether
430 U	UG/KG	Nitrobenzene	430 UJ	UG/KG	Hexachlorobenzene (HCB)
430 U -	UG/KG	Isophorone	430 U	UG/KG	Atrazine
430 U	UG/KG	2-Nitrophenol	1100 UJ	UG/KG	Pentachlorophenol
430 U	UG/KG	2,4-Dimethylphenol	430 U .	UG/KG	Phenanthrene
430 U	UG/KG	bis(2-Chloroethoxy)Methane	430 U	UG/KG	Anthracene
430 U	UG/KG	2,4-Dichlorophenol	430 U	UG/KG	Carbazole
430 U	UG/KG	Naphthalene	430 U	UG/KG	Di-n-Butylphthalate
430 U	UG/KG	4-Chloroaniline	430 U	UG/KG	Fluoranthene
430 U	UG/KG	Hexachlorobutadiene .	430 U	UG/KG	Pyrene
430 U	UG/KG	Caprolactam	430 U	UG/KG	Benzyl Butyl Phthalate
430 U	UG/KG	4-Chloro-3-Methylphenol	430 U	UG/KG	3,3'-Dichlorobenzidine
430 U	UG/KG	2-Methylnaphthalene	430 U	UG/KG	Benzo(a)Anthracene
430 UJ	UG/KG	Hexachlorocyclopentadiene (HCCP)	430 U	UG/KG	Chrysene
430 U	UG/KG	2,4,6-Trichlorophenol	430 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	430 U	UG/KG	Di-n-Octylphthalate
430 U	UG/KG	1,1-Biphenyl	430 U	UG/KG	Benzo(b)Fluoranthene
430 U	UG/KG	2-Chloronaphthalene	430 U	UG/KG	Benzo(k)Fluóranthene
1100 U	UG/KG	2-Nitroaniline	430 U	UG/KG	Benzo-a-Pyrene
430 U	UG/KĢ	Dimethyl Phthalate	430 UJ	UG/KG	Indeno (1,2,3-cd) Pyrene
430 U	UG/KĞ	2,6-Dinitrotoluene	430 UJ	UG/KG	Dibenzo(a,h)Anthracene
430 U	UG/KG	Acenaphthylene	430 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	23	%	% Moisture
	UG/KG	Acenaphthene			
	UG/KG	2,4-Dinitrophenol			
1100 ŲJ	UG/KG	4-Nitrophenol			•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6781 FY 2003

Project: 03-0474

**ANALYTE** 

STIGMAST-4-EN-3-ONE

**5 UNKNOWN COMPOUNDS** 

**MISCELLANEOUS COMPOUNDS** Facility: Gulf States Creosoting

Flowood, MS

Program: SF

190 NJ

2200 J

Id/Station: GS02SS /

RESULTS UNITS

Case No: 31635

Media: SURFACE SOIL (0" - 12")

UG/KG

UG/KG

MD No: 1XW1

D No: 1XW1

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Beginning: 04/22/2003 13:13

Requestor:

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Ending:

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample Project: 03-0474 6782 FY 2003 Requestor:

**Extractables Scan** 

Facility: Gulf States Creosoting Flowood, MS

Program: SF

Id/Station: GS02SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

Inorg Contractor: SENTIN MD No: 1XW2 D No: 1XW2

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

Project Leader: BSTRIGGO

Beginning: 04/22/2003 13:34

Ending:

DECI TO	LIMITO	ANALYTE	RESULTS	UNITS	ANALYTE
RESULTS		ANALYTE	410 U	UG/KG	Dibenzofuran
410 U	UG/KG	Benzaldehyde			
410 U	UG/KG	Phenol Phenol	410 U	UG/KG	2,4-Dinitrotoluene
410 U	UG/KG	bis(2-Chloroethyl) Ether	410 U	UG/KG	Diethyl Phthalate
410 U	UG/KG	2-Chlorophenol	410 U	UG/KG	Fluorene
410 U	UG/KG	2-Methylphenol	410 U	UG/KG	4-Chlorophenyl Phenyl Ether
410 U	UG/KG	bis(2-Chloroisopropyl) Ether	1000 UJ	UG/KG	4-Nitroaniline
410 U	UG/KG	Acetophenone	1000 U	UG/KG	2-Methyl-4,6-Dinitrophenol
410 U	UG/KG	(3-and/or 4-)Methylphenol	410 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
410 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
410 U	UG/KG	Hexachloroethane	410 U	UG/KG	4-Bromophenyl Phenyl Ether
410 U	UG/KG	Nitrobenzene	410 UJ	UG/KG	Hexachlorobenzene (HCB)
410 U.	UG/KG	Isophorone	410 U	UG/KG	Atrazine
410 U	UG/KG	2-Nitrophenol	1000 UJ	UG/KG	Pentachlorophenol
410 U	UG/KG	2,4-Dimethylphenol	410 U	UG/KG	Phenanthrene
410 U	UG/KG	bis(2-Chloroethoxy)Methane	410 U	UG/KG	Anthracene
410 U	UG/KG	2,4-Dichlorophenol	410 U	UG/KG	Carbazole
410 U	UG/KG	Naphthalene	410 U	UG/KG	Di-n-Butylphthalate
410 U	UG/KG	4-Chloroaniline	410 U	UG/KG	Fluoranthene
410 U	UG/KG	Hexachlorobutadiene	. 410 U	UG/KG	Pyrene
410 U	UG/KG	Caprolactam	410 U	UG/KG	Benzyl Butyl Phthalate
410 U	UG/KG	4-Chloro-3-Methylphenol	410 U	UG/KG	3,3'-Dichlorobenzidine
410 U	UG/KG	2-Methylnaphthalene	410 U	UG/KG	Benzo(a)Anthracene
410 UJ-	UG/KG	Hexachlorocyclopentadiene (HCCP)	410 U	UG/KG	Chrysene
410 U	UG/KG	2,4,6-Trichlorophenol	410 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1000 U	UG/KG	2,4,5-Trichlorophenol	410 U	UG/KG	Di-n-Octylphthalate
410 U	UG/KG	1,1-Biphenyl	410 U	UG/KG	Benzo(b)Fluoranthene
410 U	UG/KG	2-Chloronaphthalene	410 U	UG/KG	Benzo(k)Fluoranthene
1000 U	UG/KG	2-Nitroaniline	410 U	UG/KG	Benzo-a-Pyrene
410 U	UG/KG	Dimethyl Phthalate	410 UJ	UG/KG	Indeno (1,2,3-cd) Pyrene
410 U	UG/KG	2,6-Dinitrotoluene	410 UJ	UG/KG	Dibenzo(a,h)Anthracene
410 U	UG/KG	Acenaphthylene	410 U	UG/KG	Benzo(ghi)Perylene
1000 U	UG/KG	3-Nitroaniline	19	%	% Moisture
410 U	UG/KG	Acenaphthene	· -		•
	UG/KG	2,4-Dinitrophenol	,		
1000 UJ	UG/KG	4-Nitrophenol			
•				•	•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 13:34

Requestor:

Ending:

Sample

6782 FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

220 J

Case No: 31635

Id/Station: GS02SB /

MD No: 1XW2

Media: SUBSURFACE SOIL (> 12")

D No: 1XW2

Inorg Contractor: SENTIN Org Contractor: LIBRTY

**RESULTS UNITS** 

UG/KG

**ANALYTE** 

2 UNKNOWN OCMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 6783 FY 2003 Sample

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 14:40

Ending:

Program: SF

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

Id/Station: GS14SS /

**Extractables Scan** 

Media: SURFACE SOIL (0" - 12")

MD No: 1XW3

Inorg Contractor: SENTIN

D No: 1XW3

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
380 U	UG/KG	Benzaldehyde	380 U	UG/KG	Dibenzofuran
380 U	UG/KG	Phenol	380 U	UG/KG	2,4-Dinitrotoluene
380 U	UG/KG	bis(2-Chloroethyl) Ether	380 U	UG/KG	Diethyl Phthalate
380 U	UG/KG	2-Chlorophenol	380 U	UG/KG	Fluorene
380 U .	UG/KG	2-Methylphenol	380 U	UG/KG	4-Chlorophenyl Phenyl Ether
380 U	UG/KG	bis(2-Chloroisopropyl) Ether	970 U	UG/KG	4-Nitroaniline
380 U	UG/KG	Acetophenone	970 U	UG/KG	2-Methyl-4,6-Dinitrophenol
380 U	UG/KG	(3-and/or 4-)Methylphenol	380 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
380 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
380 U	UG/KG	Hexachloroethane	380 U	UG/KG	4-Bromophenyl Phenyl Ether
380 U	UG/KG	Nitrobenzene	380 U	UG/KG	Hexachlorobenzene (HCB)
380 U	UG/KG	Isophorone	380 U	UG/KG	Atrazine
380 U	UG/KG	2-Nitrophenol	970 U	UG/KG	Pentachlorophenol
380 U	UG/KG	2,4-Dimethylphenol	320 J	UG/KG	Phenanthrene
380 U	UG/KG	bis(2-Chloroethoxy)Methane	1100	UG/KG	Anthracene
380 U	UG/KG	2,4-Dichlorophenol	280 J	UG/KG	Carbazole
380 U	UG/KG	Naphthalene	380 U	UG/KG	Di-n-Butylphthalate
380 U	UG/KG	4-Chloroaniline	3100	UG/KG	Fluoranthene
380 U	UG/KG	Hexachlorobutadiene	3300	UG/KG	Pyrene
380 U	UG/KG	Caprolactam	380 U	UG/KG	Benzyl Butyl Phthalate
380 U	UG/KG	4-Chloro-3-Methylphenol	380 U	UG/KG	3,3'-Dichlorobenzidine
380 U	UG/KG	2-Methylnaphthalene	2300	UG/KG	Benzo(a)Anthracene
380 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	3100	UG/KG	Chrysene
380 U	UG/KG	2,4,6-Trichlorophenol	380 U	UG/KG	bis(2-Ethylhexyl) Phthalate
970 U	UG/KG	2,4,5-Trichlorophenol	380 U	UG/KG	Di-n-Octylphthalate
380 U	UG/KG	1,1-Biphenyl	4200	UG/KG	Benzo(b)Fluoranthene
380 U	UG/KG	2-Chloronaphthalene	2400	UG/KG	Benzo(k)Fluoranthene
970 U	UG/KG	2-Nitroaniline	2300	UG/KG	Benzo-a-Pyrene
380 U	UG/KG	Dimethyl Phthalate	1700	UG/KG	Indeno (1,2,3-cd) Pyrene
380 U	UG/KG	2,6-Dinitrotoluene	580	UG/KG	Dibenzo(a,h)Anthracene
380 U	UG/KG	Acenaphthylene	820	UG/KG	Benzo(ghi)Perylene
970 U	UG/KG	3-Nitroaniline	14	%	% Moisture
	UG/KG	Acenaphthene			
	UG/KG	2,4-Dinitrophenol			•
970 U	UG/KG	4-Nitrophenol			
			•		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6783 FY 2003

Facility: Gulf States Creosoting

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Flowood, MS

Program: SF

5500 J

Case No: 31635

Id/Station: GS14SS /

MD No: 1XW3

D No: 1XW3

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

UG/KG

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:40

Ending:

**RESULTS UNITS** ANALYTE 470 NJ UG/KG CYCLOPENTA (DEF) PHENANTHRENONE 1700 NJ UG/KG 11H-BENZO (A) FLUORENE 1300 NJ UG/KG PYRENE, 1-METHYL-720 NJ UG/KG PYRENE, 2-METHYL-870 NJ UG/KG 7H-BENZ [DE] ANTHRACEN-7-ONE 550 NJ UG/KG BENZO [B] NAPHTHO [2,1-D] THIOPHENE 520 NJ UG/KG 3,4-DIHYDROCYCLOPENTA (CD) PYRENE 840 NJ UG/KG TRIPHENYLENE, 2-METHYL-380 NJ UG/KG 5,12-NAPHTHACENEDIONE 2600 NJ UG/KG BENZO [E] PYRENE 710 NJ UG/KG PERYLENE 550 NJ UG/KG 1.2:7.8-DIBENZPHENANTHRENE 5600 J UG/KG **8 UNKNOWN PAHS** 

8 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

Sample 6784

FY 2003 Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 14:50

Ending:

Facility: Gulf States Creosoting Program: SF

**Extractables Scan** 

Flowood, MS

Case No: 31635

Id/Station: GS14SB /

Media: SUBSURFACE SOIL (> 12")

MD No: 1XW4

Inorg Contractor: SENTIN

D No: 1XW4

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

				<del></del>	
RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
390 U	UG/KG	Benzaldehyde	390 U	UG/KG	Dibenzofuran
390 U	UG/KG	Phenol	390 U	UG/KG	2,4-Dinitrotoluene
. 390 U	UG/KG	bis(2-Chloroethyl) Ether	390 U	UG/KG	Diethyl Phthalate
390 U	UG/KG	2-Chlorophenol	390 U	UG/KG	Fluorene
390 U	UG/KG	2-Methylphenol	390 U	UG/KG `	4-Chlorophenyl Phenyl Ether
390 U	UG/KG	bis(2-Chloroisopropyl) Ether	980 U	UG/KG	4-Nitroaniline
390 U	UG/KG	Acetophenone	980 U	UG/KG	2-Methyl-4,6-Dinitrophenol
390 U	UG/KG	(3-and/or 4-)Methylphenol	390 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
390 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	·UG/KG	1,2,4,5-Tetrachlorobenzene
390 U	UG/KG	Hexachloroethane	390 U	UG/KG	4-Bromophenyl Phenyl Ether
390 U	UG/KG	Nitrobenzene	390 U	UG/KG	Hexachlorobenzene (HCB)
. 390 U	UG/KG	Isophorone	390 U	UG/KG	Atrazine
390 U	UG/KG	2-Nitrophenol	980 U	UG/KG	Pentachlorophenol
390 U	UG/KG	2,4-Dimethylphenol	510	UG/KG	Phenanthrene
390 U *	UG/KG	bis(2-Chloroethoxy)Methane	690	UG/KG.	Anthracene
390 U	UG/KG	2,4-Dichlorophenol	210 J	UG/KG	Carbazole
390 U	UG/KG	Naphthalene	390 U	UG/KG	Di-n-Butylphthalate
· 390 U	UG/KG	4-Chloroaniline	2200	UG/KG	Fluoranthene
390 U	UG/KG	Hexachlorobutadiene	2400	UG/KG	Pyrene
390 U	UG/KG	Caprolactam	390 U	UG/KG	Benzyl Butyl Phthalate
390 U	UG/KG	4-Chloro-3-Methylphenol	390 U	UG/KG	3,3'-Dichlorobenzidine
390 U	UG/KG	2-Methylnaphthalene	1400	UG/KG	Benzo(a)Anthracene
390 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	2100	UG/KG	Chrysene
390 U	UG/KG	2,4,6-Trichlorophenol	390 U	UG/KG	bis(2-Ethylhexyl) Phthalate
980 U	UG/KG	2,4,5-Trichlorophenol	390 U	UG/KG	Di-n-Octylphthalate
390 U	UG/KG	1,1-Biphenyl	2400	UG/KG	Benzo(b)Fluoranthene
390 U	UG/KG	2-Chloronaphthalene	2000	UG/KG	Benzo(k)Fluoranthene
980 U	UG/KG	2-Nitroaniline	1400	UG/KG	Benzo-a-Pyrene
390 U	UG/KG	Dimethyl Phthalate	1100 .	UG/KG	Indeno (1,2,3-cd) Pyrene
390 U	UG/KG	2,6-Dinitrotoluene	390	UG/KG	Dibenzo(a,h)Anthracene
330 J	UG/KG	Acenaphthylene	540	UG/KG	Benzo(ghi)Perylene
980 U	UG/KG	3-Nitroaniline	15	%	% Moisture
390 U	UG/KG	Acenaphthene			
980 UJ	UG/KG	2,4-Dinitrophenol			
980 U	UG/KG	4-Nitrophenol			
		· · · · · · · · · · · · · · · · · · ·			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6784 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:50

Program: SF

Case No: 31635

Ending:

Id/Station: GS14SB /

MD No: 1XW4 D No: 1XW4.

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

MISCELLANEOUS COMPOUNDS

Org Contractor: LIBRTY

**RESULTS UNITS ANALYTE** 

320 NJ UG/KG 9,10-ANTHRACENEDIONE

320 NJ UG/KG CYCLOPENTA (dEF) PHENANTHRENONE

340 NJ PYRENE, 1-METHY-UG/KG

270 NJ **ELLIPTICINE** UG/KG

500 NJ UG/KG 7H-BENZ (DE) ANTHRACEN-7-ONE

270 NJ UG/KG BENZO (B) NAPHTHO [1,2-D] THIOPHENE 290 NJ UG/KG 3,4-DIHYDROCYCLOPENTA (CD) PYRENE

5.12-NAPHTHACENEDIONE 290 NJ UG/KG

1900 NJ UG/KG BENZO [E] PYRENE

420 NJ UG/KG PERYLENE

3200 J **7 UNKNOWN PAHS** UG/KG

340 NJ UG/KG 3900 J UG/KG

BENZO (B) TRIPHENYLENE 10 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Produced by: Goddard, Denise Sample 6785 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 15:05 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS10SS / MD No: 1XW5 Org Contractor: LIBRTY D No: 1XW5 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE		RESULTS	UNITS	ANALYTE
380 U	UG/KG	Benzaldehyde	•	120 J	UG/KG	Dibenzofuran
380 U	UG/KG	Phenol		· 380 U	UG/KG	2,4-Dinitrotoluene
380 Ū	UG/KG	bis(2-Chloroethyl) Ether		380 U	UG/KG	Diethyl Phthalate
380 U	UG/KG	2-Chlorophenol	• •	380 U	UG/KG	Fluorene
380 Ú	UG/KG	2-Methylphenol	•	380 U	UG/KG	4-Chlorophenyl Phenyl Ether
380 U	UG/KG	bis(2-Chloroisopropyl) Ether		940 U	UG/KG	4-Nitroaniline
380 U	UG/KG	Acetophenone		940 U	UG/KG	2-Methyl-4,6-Dinitrophenol
380 U	UG/KG	(3-and/or 4-)Methylphenol		380 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
380 U	UG/KG	n-Nitrosodi-n-Propylamine		NA	UG/KG	1,2,4,5-Tetrachlorobenzene
380 U	UG/KG	Hexachloroethane		380 U	UG/KG	4-Bromophenyl Phenyl Ether
380 U	UG/KG	Nitrobenzene		380 U	UG/KG	Hexachlorobenzene (HCB)
380 U	UG/KG	Isophorone		380 U	UG/KG	Atrazine
380 U	UG/KG	2-Nitrophenol	•	690 J	UG/KG	Pentachlorophenol
380 U	UG/KG	2,4-Dimethylphenol		1400	UG/KG	Phenanthrene
380 Ú	UG/KG	bis(2-Chloroethoxy)Methane		1700	UG/KG	Anthracene
380 U	UG/KG	2.4-Dichlorophenol	•	480	UG/KG	Carbazole
110 J	UG/KG	Naphthalene		380 U	UG/KG	Di-n-Butylphthalate
380 U	UG/KG	4-Chloroaniline	•	4700	UG/KG	Fluoranthene
380 U	UG/KG	Hexachlorobutadiene		5000	UG/KG	Pyrene
380 U	UG/KG	Caprolactam		380 U	UG/KG	Benzyl Butyl Phthalate
380 U	UG/KG	4-Chloro-3-Methylphenol		380 U	UG/KG	3,3'-Dichlorobenzidine
69 J	UG/KG	2-Methylnaphthalene	•	2800	UG/KG	Benzo(a)Anthracene
380 UJ	UG/KG	Hexachlorocyclopentadiene (HCCP)		4300	UG/KG	Chrysene
.380 U	UG/KG	2,4,6-Trichlorophenol		380 U	UG/KG	bis(2-Ethylhexyl) Phthalate
940 U	UG/KG	2,4,5-Trichlorophenol		380 U	UG/KG	Di-n-Octylphthalate
380 U	UG/KG	1,1-Biphenyl		5200	UG/KG	Benzo(b)Fluoranthene
380 U	UG/KG	2-Chloronaphthalene		2300	UG/KG	Benzo(k)Fluoranthene
940 U	UG/KG	2-Nitroaniline	•	3100	UG/KG	Benzo-a-Pyrene
380 U	UG/KG	Dimethyl Phthalate		2300	UG/KG	Indeno (1,2,3-cd) Pyrene
380 U	UG/KG	2,6-Dinitrotoluene	· ·	750	UG/KG	Dibenzo(a,h)Anthracene
1100	UG/KG	Acenaphthylene		1300	UG/KG	Benzo(ghi)Perylene
940 U	UG/KG	3-Nitroaniline		12	%	% Moisture
97 J	UG/KG	Acenaphthene				
940 ÚJ	UG/KG	2,4-Dinitrophenol				
940 U	UG/KG	4-Nitrophenol	•			
	UG/KG	4-Mitophenol				

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6785 FY 2003

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS10SS /

Media: SURFACE SOIL (0" - 12")

MD No: 1XW5

D No: 1XW5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 15:05

Ending:

RESULTS	UNITS	ANALYTE ,
720 J	`UG/KG	UNKNOWN CARBOXYLIC ACID
660 NJ	UG/KG	9, 10-ANTHRACENEDIONE
940 NJ	UG/KG	CYCLOPENTA (DEF) PHENANTHRENONE
2200 NJ	UG/KG	11H-BENZO [A] FLUORENE
1700 NJ	UG/KG	PYRENE, 2-METHYL-
710 NJ	UG/KG	PYRENE, 1-METHYL-
1100 NJ	UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
550 NJ	UG/KG	3,4-DIHYDROCYCLOPENTA (CD) PYRENE
800 NJ	UG/KG	TRIPHENYLENE, 2-METHYL-
560 NJ	UG/KG	2,2'-BINAPHTHALENE
690 NJ	UG/KG	5,12-NAPHTHACENEDIONE
3300 NJ	UG/KG	BENZO [E] PYRENE
1000 NJ	UG/KG	PERYLENE
4100 J	UG/KG	6 UNKNOWN COMPOUNDS
9400 J	UG/KG	10 UNKNOWN PAHS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474

Produced by: Goddard, Denise

Sample .6786 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 15:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS10SB / MD No: 1XW6 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XW6 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS** ANALYTE **RESULTS UNITS ANALYTE** 800 U UG/KG 340 J UG/KG Dibenzofuran Benzaldehyde 800 U 800 Ü UG/KG UG/KG Phenol 2.4-Dinitrotoluene Diethyl Phthalate 800 U UG/KG bis(2-Chloroethyl) Ether 800 U UG/KG 800 U 800 U UG/KG 2-Chlorophenol UG/KG Fluorene 800 U UG/KG 800 U UG/KG 4-Chlorophenyl Phenyl Ether 2-Methylphenol 800 U bis(2-Chloroisopropyl) Ether 2000 U UG/KG UG/KG 4-Nitroaniline 800 U UG/KG 2000 U UG/KG 2-Methyl-4,6-Dinitrophenol Acetophenone 800 U UG/KG 800 U UG/KG n-Nitrosodiphenylamine/Diphenylamine (3-and/or 4-)Methylphenol 800 U UG/KG 1,2,4,5-Tetrachlorobenzene UG/KG n-Nitrosodi-n-Propylamine NA 800 U UG/KG Hexachloroethane 800 U UG/KG 4-Bromophenyl Phenyl Ether 800 U UG/KG Nitrobenzene 800 U UG/KG Hexachlorobenzene (HCB) 800 U 800 U UG/KG Isophorone UG/KG Atrazine 800 U UG/KG 2-Nitrophenol 680 J UG/KG Pentachlorophenol 800 U UG/KG 2.4-Dimethylphenol 2700 UG/KG Phenanthrene 800 U UG/KG bis(2-Chloroethoxy)Methane UG/KG Anthracene 12000 800 U UG/KG 2.4-Dichlorophenol 1800 UG/KG Carbazole 390 J UG/KG Naphthalene U 008 UG/KG Di-n-Butvlphthalate 800 U UG/KG 4-Chloroaniline 28000 UG/KG Fluoranthene 800 U · UG/KG Hexachlorobutadiene UG/KG 37000 Pyrene 800 U UG/KG Caprolactam 800 U UG/KG Benzyl Butyl Phthalate 800 U UG/KG 4-Chloro-3-Methylphenol 800 U UG/KG 3,3'-Dichlorobenzidine 200 J UG/KG 2-Methylnaphthalene 23000 UG/KG Benzo(a)Anthracene 800 U UG/KG Hexachlorocyclopentadiene (HCCP) 35000 UG/KG Chrysene 800 U UG/KG 2,4,6-Trichlorophenol UG/KG 800 U bis(2-Ethylhexyl) Phthalate 2000 U UG/KG 2,4,5-Trichlorophenol 800 U UG/KG Di-n-Octvlohthalate 800 U UG/KG 1,1-Biphenyl Benzo(b)Fluoranthene 37000 UG/KG 800 U UG/KG 2-Chloronaphthalene 26000 UG/KG Benzo(k)Fluoranthene 2000 U UG/KG 2-Nitroaniline 25000 UG/KG Benzo-a-Pyrene 800 U UG/KG Dimethyl Phthalate 20000 UG/KG Indeno (1,2,3-cd) Pyrene 800 U UG/KG 2,6-Dinitrotoluene 3300 UG/KG Dibenzo(a,h)Anthracene 6100 UG/KG Acenaphthylene 5200 UG/KG Benzo(ghi)Perylene 2000 U UG/KG 3-Nitroaniline % Moisture 17 % 210 J UG/KG Acenaphthene 2000 UJ UG/KG 2,4-Dinitrophenol 2000 U UG/KG 4-Nitrophenol

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6786 FY 2003

Produced by: Goddard, Denise

MISCELLANEOUS COMPOUNDS

Requestor:

Facility: Gulf States Creosoting

Program: SF

Flowood, MS

Project Leader: BSTRIGGO

Case No: 31635

Beginning: 04/22/2003 15:15

Id/Station: GS10SB /

MD No: 1XW6

Ending:

Media: SUBSURFACE SOIL (> 12")

D No: 1XW6

Inorg Contractor: SENTIN Org Contractor: LIBRTY

RESULTS	UNITS	ANALYTE
4700 NJ	UG/KG	CYCLOPENTA (DEF) PHENANTHRENONE
9900 NJ	UG/KG	11H-BENZO [B] FLUORENE
9200 NJ	UG/KG	PYRENE, 1-METHYL-
4600 NJ	UG/KG	7H-BENZ [DE] ANTHRACENEDIONE
4600 NJ	UG/KG	BENZO [B] NAPHTHO [2,3-D] THIOPHENE
2800 NJ	UG/KG	3,4-DIHYDROCYCLOPENTA (CD) PYRENE
6000 NJ	UG/KG	CHRYSENE, 5-METHYL-
2800 NJ	UG/KG	BENZ [A] ANTHRACENE, 8-METHYL-
3100 NJ	UG/KG	2,2'-BINAPHTHALENE
4600 NJ -	UG/KG	5,12-NAPHTACENEDIONE
6200 NJ	UG/KG	BENZO [E] PYRENE
16000 NJ	UG/KG	PERYLENE
7400 NJ	UG/KG	1,2:7,8-DIBENZPHENANTHRENE
17000 J	UG/KG	4 UNKNOWN COMPOUNDS
65000 J	UG/KG	13 UNKNOWN PAHS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6787 FY 2003

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS02GW /

Media: GROUNDWATER

Case No: 31635

MD No: 1XW7 D No: 1XW7

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:55

Ending:

		<u> </u>				·		
RESULTS	UNITS	ANALYTE		·	RESULTS	UNITS	ANALYTE	
5.0 UJ	· UG/L	Benzaldehyde	•		5.0 UJ	UG/L	Dibenzofuran	
5.0 UJ	UG/L .	Phenol			5.0 UJ	UG/L	2,4-Dinitrotoluene	
5.0 UJ	UG/L	bis(2-Chloroethyl) Ether		•	5.0 UJ		Diethyl Phthalate	
5.0 UJ	UG/L	2-Chlorophenol			5.0 UJ	UG/L	Fluorene	
5.0 UJ	UG/L	2-Methylphenol			5.0 UJ		4-Chlorophenyl Phenyl Ether	
5.0 UJ	UG/L	bis(2-Chloroisopropyl) Ether	,		20 UJ		4-Nitroaniline	
5.0 UJ	UG/L	Acetophenone			20 UJ		2-Methyl-4,6-Dinitrophenol	
5.0 UJ	UG/L	(3-and/or 4-)Methylphenol			5.0 UJ		n-Nitrosodiphenylamine/Diphenylamine	
5.0 UJ	UG/L	n-Nitrosodi-n-Propylamine			5.0 UJ		1,2,4,5-Tetrachlorobenzene	
5.0 UJ	UG/L	Hexachloroethane			5.0 UJ	UG/L	4-Bromophenyl Phenyl Ether	
5.0 UJ	UG/L	Nitrobenzene	•	• .	5.0 UJ		Hexachlorobenzene (HCB)	
5.0 UJ	UG/L	Isophorone		•	5.0 UJ		Atrazine	
5.0 UJ	UG/L	2-Nitrophenol			5.0 UJ		Pentachlorophenol	•
5.0 UJ	UG/L	2,4-Dimethylphenol			5.0 UJ		Phenanthrene	·
5.0 UJ	UG/L	bis(2-Chloroethoxy)Methane	•		5.0 UJ	UG/L	Anthracene	
5.0 UJ	UG/L	2,4-Dichlorophenol			NA NA	UG/L	Carbazole '	
5.0 UJ	UG/L	Naphthalene		,	5.0 UJ	UG/L	Di-n-Butylphthalate	
5.0 UJ	UG/L	4-Chloroaniline			5.0 UJ	UG/L	Fluoranthene	
5.0 UJ	UG/L	Hexachlorobutadiene		•	5.0 UJ	UG/L	Pyrene	
5.0 UJ	UG/L	Caprolactam			5.0 UJ	UG/L	Benzyl Butyl Phthalate	
5.0 UJ	UG/L	4-Chloro-3-Methylphenol		٠	5.0 UJ	UG/L	3,3'-Dichlorobenzidine	
5.0 UJ	UG/L	2-Methylnaphthalene			5.0 UJ	UG/L	Benzo(a)Anthracene	•
5.0 UJ	UG/L	Hexachlorocyclopentadiene	(HCCP)	•	5.0 UJ	UG/L	Chrysene	
5.0 UJ	UG/L	2,4,6-Trichlorophenol	•		5.0 UJ	UG/L	bis(2-Ethylhexyl) Phthalate	
20 UJ	UG/L	2,4,5-Trichlorophenol		. *	5.0 UJ	UG/L	Di-n-Octylphthalate	
5.0 UJ	UG/L	1,1-Biphenyl	•		5.0 UJ	UG/L	Benzo(b)Fluoranthene	
5.0 UJ	UG/L	2-Chloronaphthalene	·		5.0 UJ	UG/L	Benzo(k)Fluoranthene	
20 UJ	UG/L	2-Nitroaniline			5.0 UJ	UG/L	Benzo-a-Pyrene	
5.0 UJ `	UG/L	Dimethyl Phthalate			5.0 UJ	UG/L	Indeno (1,2,3-cd) Pyrene	
5.0 UJ	UG/L	2,6-Dinitrotoluene			5.0 UJ	UG/L	Dibenzo(a,h)Anthracene	•
5.0 UJ	UG/L	Acenaphthylene			5.0 UJ	UG/L	Benzo(ghi)Perylene	
20 UJ	UG/L	3-Nitroaniline					1.5 / 1	
5.0 UJ	UG/L	Acenaphthene					•	
20 UJ	UG/L	2,4-Dinitrophenol					•	
20 UJ	UG/L	4-Nitrophenol						
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U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Project: 03-0474

6787 FY 2003

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS02GW /

Media: GROUNDWATER

Flowood, MS

Case No: 31635 MD No: 1XW7

D No: 1XW7 ·

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:55

Ending:

**RESULTS UNITS ANALYTE** 2.2 NJ UG/L LIMONENE

Data Reported as Identified by CLP Lab - IDs Not Verified

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L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6788 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:00 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS11SS / MD No: 1XW8 Org Contractor: LIBRTY D No: 1XW8 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
350 U	UG/KG	Benzaldehyde	140 J	UG/KG	Dibenzofuran
350 U	UG/KG	Phenol	350 U	UG/KG	2,4-Dinitrotoluene
350 U	UG/KG	bis(2-Chloroethyl) Ether	350 U	UG/KG	Diethyl Phthalate
350 U	UG/KG	2-Chlorophenol	350 U	UG/KG	Fluorene
350 U	UG/KG	2-Methylphenol	350 U	UG/KG	4-Chlorophenyl Phenyl Ether
350 U	UG/KG	bis(2-Chloroisopropyl) Ether	890 U	UG/KG	4-Nitroaniline
350 U	`UG/KG	Acetophenone	890 U	UG/KG	2-Methyl-4,6-Dinitrophenol
350 U	UG/KG	(3-and/or 4-)Methylphenol	350 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
350 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
350 U	UG/KG	Hexachloroethane	350 U	UG/KG	4-Bromophenyl Phenyl Ether
350 U `	UG/KG	Nitrobenzene	350 U	UG/KG	Hexachlorobenzene (HCB)
350 U	UG/KG	Isophorone	350 U	UG/KG	Atrazine
350 U	UG/KG	2-Nitrophenol	640 J	UG/KG	Pentachlorophenol
350 U	UG/KG	2,4-Dimethylphenol	540	UG/KG	Phenanthrene
350 U	ÚG/KG	bis(2-Chloroethoxy)Methane	830	.UG/KG	Anthracene
350 U	UG/KG	2,4-Dichlorophenol	210 J	UG/KG	Carbazole
120 J	UG/KG	Naphthalene	350 U	UG/KG	Di-n-Butylphthalate
350 U	UG/KG	4-Chloroaniline	1900	UG/KG	Fluoranthene
350 U	UG/KG	Hexachlorobutadiene	1700	UG/KG	Pyrene
350 U	UG/KG	Caprolactam	350 U	UG/KG	Benzyl Butyl Phthalate
350 U	UG/KG	4-Chloro-3-Methylphenol	350 U	UG/KG	3,3'-Dichlorobenzidine
- 70 J	UG/KG	2-Methylnaphthalene	1000	UG/KG	Benzo(a)Anthracene
350 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	1500	UG/KG	Chrysene
350 U	UG/KG	2,4,6-Trichlorophenol	350 U	UG/KG	bis(2-Ethylhexyl) Phthalate
890 U	UG/KG	2,4,5-Trichlorophenol	350 U	UG/KG	Di-n-Octylphthalate
350 U	UG/KG	1,1-Biphenyl	2500	UG/KG	Benzo(b)Fluoranthene
350 U	UG/KG	2-Chloronaphthalene	1700	UG/KG	Benzo(k)Fluoranthene
890 U	UG/KG	2-Nitroaniline	800	UG/KG	Benzo-a-Pyrene
350 U	UG/KG	Dimethyl Phthalate	940	UG/KG	Indeno (1,2,3-cd) Pyrene
350 U	UG/KG	2,6-Dinitrotoluene	310 J	UG/KG	Dibenzo(a,h)Anthracene
330 J	UG/KG	Acenaphthylene	460	·UG/KG	Benzo(ghi)Perylene
890 U	UG/KG	3-Nitroaniline	7	%	% Moisture
350 U	UG/KG	Acenaphthene			
	UG/KG	2,4-Dinitrophenol			
890 U	UG/KG	4-Nitrophenol			
					•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

Sample 6788 FY 2003 Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

3600 J

Id/Station: GS11SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

UG/KG

MD No: 1XW8 D No: 1XW8

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:00

Ending:

RESULTS	UNITS	ANALYTE
670 NJ	UG/KG	HEXADECANOIC ACID
260 NJ *	UG/KG	9, 10-ANTHRACENEDIONE
320 NJ	UG/KG	CYCLOPENTA (DEF) PHENANTHRENONE
390 NJ	UG/KG	PYRENE, 1-METHYL-
430 NJ	UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
370 NJ	UG/KG	TRIPHENYLENE, 2-METHYL-
330 NJ	UG/KG	BENZ (A) ANTHRACENE-7,12-DIONE
1700 NJ	UG/KG	BENZO [E] PYRENE
480 J	UG/KG	1,2:7,8-DIBENZPHENANTHRRENE
380 NJ	UG/KG	BENZO [B] TRIPHENYLENE
2100 J	UG/KG	6 UNKNOWN PAHS

12 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

400 U

400 U

400 U

400 U

400 U

400 U

400 U

400 U

1000 U

400 U

400 U

400 U

400 U

1000 UJ

1000 U

1000 U

1000 U

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

Hexachlorobutadiene

2-Methylnaphthalene

2,4,6-Trichlorophenol

2,4,5-Trichlorophenol

2-Chloronaphthalene

Dimethyl Phthalate

2.6-Dinitrotoluene

Acenaphthylene

2,4-Dinitrophenol

4-Chioro-3-Methylphenol

Hexachlorocyclopentadiene (HCCP)

Caprolactam

1.1-Biphenyl

2-Nitroaniline

3-Nitroaniline

Acenaphthene

4-Nitrophenol

Produced by: Goddard, Denise 6789 FY 2003 Sample Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:10 Program: SF Case No: 31635 Ending: Id/Station: GS11SB / Inorg Contractor: SENTIN MD No: 1XW9 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XW9 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE RESULTS UNITS ANALYTE** 400 Ú UG/KG Benzaldehyde 400 U UG/KG Dibenzofuran 400 U UG/KG 400 U UG/KG 2.4-Dinitrotoluene Phenol 400 U UG/KG bis(2-Chloroethyl) Ether 400 U UG/KG Diethyl Phthalate 400 U 400 U UG/KG Fluorene UG/KG 2-Chlorophenol 400 U UG/KG 4-Chlorophenyl Phenyl Ether 400 U UG/KG 2-Methylphenol 400 U bis(2-Chloroisopropyl) Ether 1000 U UG/KG 4-Nitroaniline UG/KG 400 U 1000 U UG/KG 2-Methyl-4,6-Dinitrophenol UG/KG Acetophenone 400 U UG/KG (3-and/or 4-)Methylphenol 400 U UG/KG n-Nitrosodiphenvlamine/Diphenvlamine UG/KG 1.2.4.5-Tetrachlorobenzene 400 U UG/KG n-Nitrosodi-n-Propylamine NA 400 U UG/KG 4-Bromophenyl Phenyl Ether 400 U UG/KG Hexachioroethane 400 U UG/KG Nitrobenzene 400 U UG/KG Hexachlorobenzene (HCB) 400 U 400 U UG/KG Atrazine UG/KG Isophorone Pentachlorophenol 400 U UG/KG 2-Nitrophenol 1000 U UG/KG 400 U 2,4-Dimethylphenol 400 U UG/KG UG/KG Phenanthrene 400 U UG/KG bis(2-Chloroethoxy)Methane 49 J UG/KG Anthracene 400 U UG/KG 2,4-Dichlorophenol 400 U UG/KG Carbazole 400 U UG/KG Naphthalene 400 U UG/KG Di-n-Butylphthalate 400 U UG/KG 4-Chloroaniline 400 U UG/KG Fluoranthene

400 U

400 U

400 U

400 U

400 U

400 U

400 U

400 U

46 J

400 U

400 U

400 U

400 U

17

UG/KG

UG/KG

UG/KG

UG/KG

.UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

%

Pvrene

Chrysene

Benzyl Butyl Phthalate

3,3'-Dichlorobenzidine

bis(2-Ethylhexyl) Phthalate

Benzo(a)Anthracene

Di-n-Octylphthalate

Benzo-a-Pyrene

% Moisture

Benzo(b)Fluoranthene

Benzo(k)Fluoranthene

Indeno (1,2,3-cd) Pyrene

Dibenzo(a,h)Anthracene

Benzo(ghi)Perylene

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

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**MISCELLANEOUS COMPOUNDS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6789

FY **2003** 

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:10

Ending:

Id/Station: GS11SB /

Program: SF

530 J

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

MD No: 1XW9

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

UG/KG

D No: 1XW9

Org Contractor: LIBRTY

**RESULTS UNITS** 86 J UG/KG

**ANALYTE** 

UNKNOWN CARBOXYLIC ACID 3 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample 6790 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 16:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS09SS / MD No: 1XX0 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX0 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS** ANALYTE **RESULTS UNITS** ANALYTE 450 U UG/KG Benzaldehyde 450 U UG/KG Dibenzofuran 450 Ù UG/KG 2.4-Dinitrotoluene 450 U - UG/KG Phenol 450 U UG/KG .Diethyl Phthalate 450 U UG/KG bis(2-Chloroethyl) Ether 450 U UG/KG 2-Chlorophenol 450 U UG/KG Fluorene 450 U UG/KG 4-Chlorophenyl Phenyl Ether 450 U UG/KG 2-Methylphenol bis(2-Chloroisopropyl) Ether 1100 U UG/KG 4-Nitroaniline 450 U UG/KG 1100 U UG/KG 2-Methyl-4,6-Dinitrophenol 450 U UG/KG Acetophenone n-Nitrosodiphenylamine/Diphenylamine 450 U UG/KG 450 U UG/KG (3-and/or 4-)Methylphenol 450 U UG/KG n-Nitrosodi-n-Propylamine NA UG/KG 1.2.4.5-Tetrachlorobenzene UG/KG 4-Bromophenyl Phenyl Ether 450 U UG/KG Hexachloroethane 450 U 450 U 450 U UG/KG Hexachlorobenzene (HCB) UG/KG Nitrobenzene 450 U UG/KG Atrazine 450 U UG/KG Isophorone 1100 U UG/KG 450 U 2-Nitrophenol Pentachlorophenol UG/KG 450 U UG/KG 2,4-Dimethylphenol 450 U UG/KG Phenanthrene 450 U bis(2-Chloroethoxy)Methane 450 U UG/KG UG/KG Anthracene 450 U UG/KG 2,4-Dichlorophenol 450 U UG/KG Carbazole 450 U 450 U UG/KG Naphthalene UG/KG Di-n-Butylphthalate 450 U UG/KG 4-Chloroaniline 450 U UG/KG Fluoranthene 450 U UG/KG Hexachlorobutadiene 450 U UG/KG Pyrene 450 U 450 U UG/KG Benzyl Butyl Phthalate UG/KG Caprolactam 450 U 450 U UG/KG 4-Chloro-3-Methylphenol UG/KG 3,3'-Dichlorobenzidine 450 U UG/KG 2-Methylnaphthalene 450 U UG/KG Benzo(a)Anthracene 450 U UG/KG Hexachlorocyclopentadiene (HCCP) 450 U UG/KG Chrysene 450 U UG/KG 2,4,6-Trichlorophenol 450 U UG/KG bis(2-Ethylhexyl) Phthalate 1100 U 450 U UG/KG 2,4,5-Trichlorophenol UG/KG Di-n-Octylphthalate 450 U UG/KG 450 U UG/KG Benzo(b)Fluoranthene 1,1-Biphenyl 450 U 2-Chloronaphthalene UG/KG 450 U UG/KG Benzo(k)Fluoranthene 1100 U UG/KG 2-Nitroaniline 450 U UG/KG Benzo-a-Pyrene 450 U UG/KG Dimethyl Phthalate 450 U UG/KG Indeno (1,2,3-cd) Pyrene 450 U 2,6-Dinitrotoluene UG/KG 450 U UG/KG Dibenzo(a,h)Anthracene 450 U UG/KG 450 U Acenaphthylene UG/KG Benzo(ghi)Perylene 1100 U UG/KG 3-Nitroaniline 26 % % Moisture 450 U UG/KG Acenaphthene 1100 UJ 2,4-Dinitrophenol UG/KG 1100 U UG/KG 4-Nitrophenol

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

EPA - REGION IV SESD, ATHENS, GA

Production Date: 06/20/2003 14:10

Sample

6790 FY 2003

Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS09SS /

Media: SURFACE SOIL (0" - 12")

Case No: 31635 MD No: 1XX0

D No: 1XX0

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:15

Ending:

**RESULTS UNITS** ANALYTE

97 NJ . UG/KG PENTADECANOIC ACID

140 NJ UG/KG 1-TETRADECENE 1-OCTADECENE 230 NJ UG/KG 230 NJ UG/KG 1-HEPTADECENE

1900 J UG/KG

10 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Facility: Gulf States Creosoting

Program: SF

Produced by: Goddard, Denise Sample 6791 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** 

Project Leader: BSTRIGGO Flowood, MS Beginning: 04/22/2003 16:25 Case No: 31635

Ending:

Id/Station: GS09SB / Inorg Contractor: SENTIN MD No: 1XX1

Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XX1 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
420 U	UG/KG	Benzaldehyde	420 U	UG/KG	Dibenzofuran
420 U	UG/KG	Phenol	420 U	UG/KG	2,4-Dinitrotoluene
420 U	UG/KG	bis(2-Chloroethyl) Ether	420 U	UG/KG	Diethyl Phthalate
420 U	UG/KG	2-Chlorophenol	420 U	UG/KG	Fluorene
420 U	UG/KG	2-Methylphenol	420 U	UG/KG	4-Chlorophenyl Phenyl Ether
420 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
420 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
420 U	UG/KG	(3-and/or 4-)Methylphenol	420 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
420 U	UG/KG	n-Nitrosodi-n-Propylamine	NA.	UG/KG	1,2,4,5-Tetrachlorobenzene
420 U	UG/KG	Hexachloroethane	420 U	UG/KG	4-Bromophenyl Phenyl Ether
420 U	UG/KG	Nitrobenzene	420 U	UG/KG	Hexachlorobenzene (HCB)
420 U	· UG/KG	Isophorone	420 U	UG/KG	Atrazine
420 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
420 Ū	UG/KG	2,4-Dimethylphenol	420 U	UG/KG	Phenanthrene
420 U	UG/KG	bis(2-Chloroethoxy)Methane	420 U	UG/KG	Anthracene
420 U	UG/KG	2,4-Dichlorophenol	420 U	UG/KG	Carbazole
420 U	UG/KG	Naphthalene	420 U	UG/KG	Di-n-Butylphthalate
420 U	UG/KG	4-Chloroaniline	420 U	UG/KG	Fluoranthene
420 U	UG/KG	Hexachlorobutadiene	420 U	UG/KG	Pyrene
420 U	UG/KG	Caprolactam	420 U	UG/KG	Benzyl Butyl Phthalate
420 U	UG/KG	4-Chloro-3-Methylphenol	420 U	UG/KG	3,3'-Dichlorobenzidine
420 U	UG/KG	2-Methylnaphthalene	420 U	UG/KG	Benzo(a)Anthracene
420 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	420 U	UG/KG	Chrysene
420 U	UG/KG	2,4,6-Trichlorophenol	420 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	420 U	UG/KG	Di-n-Octylphthalate
420 U	UG/KG	1,1-Biphenyl	420 U	UG/KG	Benzo(b)Fluoranthene
420 U	UG/KG	2-Chloronaphthalene	420 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	420 U	UG/KG	Benzo-a-Pyrene
420 U	UG/KG	Dimethyl Phthalate	420 U	UG/KG	Indeno (1,2,3-cd) Pyrene
420 U	UG/KG	2,6-Dinitrotoluene	420 U	UG/KG	Dibenzo(a,h)Anthracene
420 U	UG/KG	Acenaphthylene	420 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	21	%	% Moisture
420 U	UG/KG	Acenaphthene			
1100 UJ	UG/KG	2,4-Dinitrophenol	•	_	·
1100 U	UG/KG	4-Nitrophenol			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6791 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:25

Ending:

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

ld/Station: GS09SB /

MD No: 1XX1

Media: SUBSURFACE SOIL (> 12")

D No: 1XX1

Inorg Contractor: SENTIN Org Contractor: LIBRTY

**RESULTS UNITS** 

**ANALYTE** 

370 J UG/KG 3 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6792 FY 2003 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 15:15 Program: SF Case No: 31635 Ending: Id/Station: GS06SD / Inorg Contractor: SENTIN MD No: 1XX2

Media: SEDIMENT D No: 1XX2 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

330 U G/KG   Benzaldehyde   330 U UG/KG   Dibenzofuran   330 U UG/KG   Didenzofuran   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   2.4-Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   3.2-Mithyl-1.4-Dintrotoluene   330 U UG/KG   3.3-Mitrosoluene   330 U UG/KG   3.3-Mitrosoluene   330 U UG/KG   3.3-Mitrosoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   330 U UG/KG   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintrotoluene   4Dintro	RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
330 U   UG/KG   Diethyl Phthalate   330 U   UG/KG   Chloropenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophenol   330 U   UG/KG   Chlorophe	330 U	UG/KG	Benzaldehyde	330 U	UG/KG	Dibenzofuran
330 U   UG/KG   2-Chlorophenol   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   3-Chlorophenol   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   4-Chlorophenol   330 U   UG/KG   3-Chlorophenol   330 U   UG/KG	330 U	UG/KG	Phenol			
1	330 U	UG/KG	bis(2-Chloroethyl) Ether	330 U		Diethyl Phthalate
330 U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone   San U   UG/KG   Acetophenone	330 U	UG/KG	2-Chlorophenol		UG/KG	Fluorene
330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-)Methylphenol   330 U   UG/KG   Cand/or 4-Methylphenol   330	330 U	UG/KG	2-Methylphenol		UG/KG	4-Chlorophenyl Phenyl Ether
330 U   UG/KG   1-Nitrosodin-n-Propylamine   330 U   UG/KG   1-2,4,5-Tetrachlorobenzene   330 U   UG/KG   1-2,4,5-Tetrachlorobenzene   330 U   UG/KG   1-2,4,5-Tetrachlorobenzene   330 U   UG/KG   1-2,4,5-Tetrachlorobenzene   330 U   UG/KG   1-2,4,5-Tetrachlorobenzene   330 U   UG/KG   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlorobenzene   1-2,4,5-Tetrachlor	330 U	UG/KG	bis(2-Chloroisopropyl) Ether			4-Nitroaniline
330 U   UG/KG   Hexachloroethane   330 U   UG/KG   Hexachloroethane   330 U   UG/KG   Hexachloroethane   330 U   UG/KG   Hexachloroethane   330 U   UG/KG   Hexachloroethane   330 U   UG/KG   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloroethane   Hexachloro	330 U	UG/KG	Acetophenone		UG/KG	2-Methyl-4,6-Dinitrophenol
330 U   UG/KG   Hexachloroethane   330 U   UG/KG   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexachloropena   Hexa		UG/KG				
330 U   UG/KG   Senzo   UG/KG   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senzo   Senz		UG/KG	n-Nitrosodi-n-Propylamine			
330 U		UG/KG	Hexachloroethane			4-Bromophenyl Phenyl Ether
330 U   UG/KG   2-Nitrophenol   25 U   UG/KG   Pentachlorophenol   Phenanthrene   330 U   UG/KG   Phenanthrene   330 U   UG/KG   Phenanthrene   330 U   UG/KG   Phenanthrene   330 U   UG/KG   Phenanthrene   330 U   UG/KG   Phenanthrene   330 U   UG/KG   Phenanthrene   Anthracene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanthrene   Phenanth		UG/KG	Nitrobenzene			Hexachlorobenzene (HCB)
330 U   UG/KG   2,4-Dimethylphenol   330 U   UG/KG   Anthracene   330 U   UG/KG   Anthracene   330 U   UG/KG   Anthracene   330 U   UG/KG   Carbazole   330 U   UG/KG   Carbazole   330 U   UG/KG   Carbazole   330 U   UG/KG   Din-Butylphthalate   330 U   UG/KG   Din-Butylphthalate   330 U   UG/KG   Din-Butylphthalate   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   330 U   UG/KG   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene   Prene						
330 U   UG/KG   Dis(2-Chloropthoxy)Methane   330 U   UG/KG   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Carbazole   Car						
330 U UG/KG						
330 U   UG/KG   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable   Vacable						
330 U         UG/KG         4-Chloroaniline         330 U         UG/KG         Fluoranthene           330 U         UG/KG         Hexachlorobutadiene         330 U         UG/KG         Pyrene           330 U         UG/KG         Caprolactam         330 U         UG/KG         Benzyl Butyl Phthalate           330 U         UG/KG         4-Chloro-3-Methylphenol         330 U         UG/KG         Benzo(a)Anthracene           330 U         UG/KG         2-Methylnaphthalene         330 U         UG/KG         Benzo(a)Anthracene           330 U         UG/KG         2-4,6-Trichlorophenol         330 U         UG/KG         Chrysene           330 U         UG/KG         2,4,5-Trichlorophenol         330 U         UG/KG         Di-n-Octylphthalate           330 U         UG/KG         1,1-Biphenyl         330 U         UG/KG         Benzo(b)Fluoranthene           330 U         UG/KG         2-Chloronaphthalene         330 U         UG/KG         Benzo(k)Fluoranthene           830 U         UG/KG         2-Nitroaniline         330 U         UG/KG         Benzo-a-Pyrene           330 U         UG/KG         2,6-Dinitrotoluene         330 U         UG/KG         Dibenzo(a,h)Anthracene           330 U         UG/KG						
330 U   UG/KG   Hexachlorobutadiene   330 U   UG/KG   Caprolactam   330 U   UG/KG   Benzyl Butyl Phthalate   330 U   UG/KG   Hoxachloro-3-Methylpheno    330 U   UG/KG   Hoxachloro-2-Methylpheno    330 U   UG/KG   Hoxachloro-2Methylpheno    330 U   UG/KG   Hoxachloro-2Methylpheno    330 U   UG/KG   Hoxachloro-2Methylpheno    330 U   UG/KG   Hoxachloro-2						
330 U UG/KG Caprolactam 330 U UG/KG 4-Chloro-3-Methylphenol 330 U UG/KG 2-Methylnaphthalene 330 U UG/KG 2-Methylnaphthalene 330 U UG/KG Hexachlorocyclopentadiene (HCCP) 330 U UG/KG 2,4,6-Trichlorophenol 330 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG 1,1-Biphenyl 330 U UG/KG 2-Chloronaphthalene 330 U UG/KG 2-Chloronaphthalene 330 U UG/KG 2-Nitroaniline 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG 3-Nitroaniline 330 U UG/KG Acenaphthylene 330 U UG/KG Acenaphthylene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG 2,4-Dinitrophenol						
330 U UG/KG 2-Methylnaphthalene 330 U UG/KG Benzo(a)Anthracene 330 U UG/KG Hexachlorocyclopentadiene (HCCP) 330 U UG/KG Chrysene 330 U UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/KG UG/			:	330 U		
330 U UG/KG 2-Methylnaphthalene 330 U UG/KG Benzo(a)Anthracene 330 U UG/KG Chrysene 330 U UG/KG Chrysene 330 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG Di-n-Octylphthalate 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Benzo(a)Anthracene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(a)Anthracene 330 U UG/KG Benzo(a)Anthracene 330 U UG/KG Benzo(a)Anthracene 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Acenaphthylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Acenaphthene 330 U UG/KG 3-Nitroaniline 20 % Moisture 330 U UG/KG 2,4-Dinitrophenol						
330 U UG/KG Hexachlorocyclopentadiene (HCCP) 330 U UG/KG 2,4,6-Trichlorophenol 330 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG Din-Octylphthalate 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(k)Fluoranthene 330 U UG/KG Benzo(k)Fluoranthene 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Acenaphthylene 330 U UG/KG Acenaphthene 330 U UG/KG Z,4-Dinitrophenol						
330 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG Di-n-Octylphthalate 330 U UG/KG 1,1-Biphenyl 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG 2-Chloronaphthalene 330 U UG/KG Benzo(k)Fluoranthene 330 U UG/KG 2-Nitroaniline 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Indeno (1,2,3-cd) Pyrene 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Acenaphthylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Acenaphthene 330 U UG/KG 3-Nitroaniline 20 % Moisture						
830 U UG/KG 2,4,5-Trichlorophenol 330 U UG/KG Di-n-Octylphthalate 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG Benzo(c)Fluoranthene 330 U UG/KG Benzo(c)Fluoranthene 330 U UG/KG Benzo(a,b)Fluoranthene 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Indeno (1,2,3-cd) Pyrene 330 U UG/KG Dibenzo(a,b)Anthracene 330 U UG/KG Dibenzo(a,b)Anthracene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene						
330 U UG/KG 1,1-Biphenyl 330 U UG/KG Benzo(b)Fluoranthene 330 U UG/KG 2-Chloronaphthalene 830 U UG/KG 2-Nitroaniline 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Indeno (1,2,3-cd) Pyrene 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Acenaphthylene 330 U UG/KG Benzo-(a,h)Anthracene 330 U UG/KG Benzo(a,h)Anthracene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene						
330 U UG/KG 2-Chloronaphthalene 830 U UG/KG 2-Nitroaniline 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dimethyl Phthalate 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG Acenaphthylene 330 U UG/KG 3-Nitroaniline 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG 2,4-Dinitrophenol						
830 U UG/KG 2-Nitroaniline 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG Dimethyl Phthalate 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG Acenaphthylene 330 U UG/KG 3-Nitroaniline 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Benzo-a-Pyrene 330 U UG/KG Dibenzo(a,h)Anthracene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG 2,4-Dinitrophenol						
330 U UG/KG Dimethyl Phthalate 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG Acenaphthylene 330 U UG/KG Acenaphthylene 330 U UG/KG 3-Nitroaniline 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene						
330 U UG/KG 2,6-Dinitrotoluene 330 U UG/KG Acenaphthylene 330 U UG/KG Acenaphthylene 330 U UG/KG Benzo(ghi)Perylene 330 U UG/KG 3-Nitroaniline 330 U UG/KG Acenaphthene 330 U UG/KG Acenaphthene 330 U UG/KG 2,4-Dinitrophenol						
330 U UG/KG Acenaphthylene  830 U UG/KG 3-Nitroaniline  330 U UG/KG Benzo(ghi)Perylene  20 % Moisture  330 U UG/KG Acenaphthene  830 UJ UG/KG 2,4-Dinitrophenol						
830 U UG/KG 3-Nitroaniline 20 % Moisture 330 U UG/KG Acenaphthene 830 UJ UG/KG 2,4-Dinitrophenol			,			
330 U ÚG/KG Acenaphthene 830 UJ ÚG/KG 2,4-Dinitrophenol						
830 UJ UG/KG 2,4-Dinitrophenol				20	%	% Moisture
					÷	
osu u ug/kg 4-ivitropnenoi						•
	830 0	UG/KG	4-Nitropnenor			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Project: 03-0474

6792 FY 2003

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS06SD /

MD No: 1XX2

Media: SEDIMENT

D No: 1XX2

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Beginning: 04/22/2003 15:15 Ending:

Requestor:

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

**RESULTS UNITS** 61 J

UG/KG

**ANALYTE** 

**UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

Sample

6792 FY 2003

Project: 03-0474

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS06SD / Media: SEDIMENT

**SPECIFIED TESTS** 

Case No: 31635 MD No: 1XX2

D No: 1XX2

Inorg Contractor: SENTIN Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 15:15

Requestor:

Ending:

RESULTS	UNITS	ANALYTE
.0.4 J	UG/KG	2-Methylnaphthalene
0.5 J	UG/KG	Naphthalene
0.8 U	UG/KG	Acenaphthylene
0.8 U	UG/KG	Acenaphthene
0.3 J	UG/KG	Fluorene
4	UG/KG	Phenanthrene
. 2 U	UG/KG	Anthracene
12 U	UG/KG	Fluoranthene
10 U	UG/KG	Pyrene
5 U	UG/KG	Benzo(a)Anthracene
7 U	UG/KG	Chrysene
NA	UG/KG	Benzo(b)Fluoranthene
NA	UG/KG	Benzo(k)Fluoranthene
4 U	UG/KG	Benzo-a-Pyrene
NA	UG/KG	Indeno (1,2,3-cd) Pyrene
1	UG/KG	Dibenzo(a,h)Anthracene
,NA	UG/KG	Benzo(ghi)Perylene
13 U	UG/KG	Pentachlorophenol
		· ·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample 6793 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 15:50 Program: SF Case No: 31635 Ending: Id/Station: GS08SD / MD No: 1XX3 Inorg Contractor: SENTIN Org Contractor: LIBRTY Media: SEDIMENT D No: 1XX3 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
330 U	UG/KG	Benzaldehyde	330 U.	UG/KG	Dibenzofuran
330 U	UG/KG	Phenol	330 U	UG/KG	2.4-Dinitrotoluene
330 U	UG/KG	bis(2-Chloroethyl) Ether	330 U	UG/KG	Diethyl Phthalate
330 U	.UG/KG	2-Chlorophenol	330 U	UG/KG	Fluorene
330 U	UG/KG	2-Methylphenol	330 U	UG/KG	4-Chlorophenyl Phenyl Ether
330 U	UG/KG	bis(2-Chloroisopropyl) Ether	830 U	UG/KG	4-Nitroaniline
330 U	UG/KG	Acetophenone	830 U	UG/KG	2-Methyl-4,6-Dinitrophenol
330 U	UG/KG	(3-and/or 4-)Methylphenol	330 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
330 Ū	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
330 U	UG/KG	Hexachloroethane	330 U	UG/KG	4-Bromophenyl Phenyl Ether
330 U	UG/KG	Nitrobenzene	330 U	UG/KG	Hexachlorobenzene (HCB)
330 U	UG/KG	Isophorone	330 U	UG/KG	Atrazine
330 Ü	UG/KG	2-Nitrophenol	830 U	UG/KG	Pentachlorophenol
330 U	UG/KG	2,4-Dimethylphenol	330 U	UG/KG	Phenanthrene
330 U	UG/KG	bis(2-Chloroethoxy)Methane	330 U	UG/KG	Anthracene
330 U	UG/KG	2,4-Dichlorophenol	330 U	UG/KG	Carbazole
330 U	UG/KG	Naphthalene	330 U	UG/KG	Di-n-Butylphthalate
330 U	UG/KG	4-Chloroaniline	330 Ū	UG/KG	Fluoranthene
330 U	UG/KG	Hexachlorobutadiene	330 U	UG/KG	Pyrene
- 330 U	UG/KG	Caprolactam	330 U	UG/KG	Benzyl Butyl Phthalate
330 U	UG/KG	4-Chloro-3-Methylphenol	330 U	UG/KG	3,3'-Dichlorobenzidine
330 U .	UG/KG	2-Methylnaphthalene	330 Ū	UG/KG	Benzo(a)Anthracene
330 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	330 U	UG/KG	Chrysene
330 U	UG/KG	2,4,6-Trichlorophenol	330 Ū	UG/KĠ	bis(2-Ethylhexyl) Phthalate
830 U	UG/KG	2,4,5-Trichlorophenol	330 U	UG/KG	Di-n-Octylphthalate
330 U	UG/KG	1,1-Biphenyl	330 U	UG/KG	Benzo(b)Fluoranthene
330 U	UG/KG	2-Chloronaphthalene	330 U	UG/KG	Benzo(k)Fluoranthene
830 U	UG/KG	2-Nitroaniline	330 U	UG/KG	Benzo-a-Pyrene
330 U	UG/KG	Dimethyl Phthalate	330 U	UG/KG	Indeno (1,2,3-cd) Pyrene
330 U	UG/KG	2.6-Dinitrotoluene	330 U	UG/KG	Dibenzo(a,h)Anthracene
330 U	UG/KG	Acenaphthylene	330 U	UG/KG	Benzo(ghi)Perylene
830 U	UG/KG	3-Nitroaniline	19	%	% Moisture
330 U	UG/KG	Acenaphthene			,
830 UJ	UG/KG	2,4-Dinitrophenol			
830 U	UG/KG	4-Nitrophenol			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6793 FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS08SD / Media: SEDIMENT

MD No: 1XX3

D No: 1XX3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 15:50

Ending:

**RESULTS UNITS** 

170 J

**ANALYTE** UG/KG

3 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not defected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 06/20/2003 14:10

Sample **SPECIFIED TESTS** 

6793 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 15:50

Ending:

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

Program: SF Id/Station: GS08SD /

Media: SEDIMENT

Inorg Contractor: SENTIN MD No: 1XX3

Org Contractor: LIBRTY D No: 1XX3

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
0.8 U	UG/KG	2-Methylnaphthalene
0.4 J	UG/KG	Naphthalene
0.8 U	UG/KG	Acenaphthylene
0.8 U	· UG/KG	Acenaphthene
0.8 U	UG/KG	Fluorene
0.0 U	UG/KG	Phenanthrene
0.8 U	UG/KG	Anthracene
2 U	UG/KG	Fluoranthene
2 U	UG/KG	Pyrene
1 U	UG/KG	Benzo(a)Anthracene
2 U	UG/KG	Chrysene
NA	UG/KG	Benzo(b)Fluoranthene
NA	UG/KG	Benzo(k)Fluoranthene
0.8 U	UG/KG	Benzo-a-Pyrene
NA	UG/KG	Indeno (1,2,3-cd) Pyrene
0.8 U	UG/KG	Dibenzo(a,h)Anthracene
NA	UG/KG	Benzo(ghi)Perylene
13 U	UG/KG	Pentachlorophenol

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Project: 03-0474 Sample FY **2003** Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:42 Program: SF Case No: 31635 Ending: Id/Station: GS22SS / Inorg Contractor: SENTIN MD No: 1XX4

Media: SURFACE SOIL (0" - 12") D No: 1XX4 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	<del></del>	· · · · · · · · · · · · · · · · · · ·	· <del></del>	RESULTS	UNITS	ANALYTE	
380 U	UG/KG	Benzaldehyde		• • •	•	380 U	UG/KG	Dibenzofuran	
380 U	UG/KG	Phenol			•	380 U	UG/KG	2,4-Dinitrotoluene	
380 U	UG/KG	bis(2-Chloroethyl) Ether	İ			380 U	UG/KG	Diethyl Phthalate	
380 U	UG/KG	2-Chlorophenol				380 U	UG/KG	Fluorene	•
380 U	UG/KG	2-Methylphenol		•		380 U	UG/KG	4-Chlorophenyl Phenyl Ether	
380 U	UG/KG	bis(2-Chloroisopropyl) E	thor			950 U	UG/KG	4-Nitroaniline	
380 U	UG/KG UG/KG	Acetophenone	riei			950 U	UG/KG	2-Methyl-4,6-Dinitrophenol	
380 U	UG/KG UG/KG					380 U	UG/KG		
		(3-and/or 4-)Methylphen				NA	UG/KG	n-Nitrosodiphenylamine/Diphenylamine	
380 U	UG/KG	n-Nitrosodi-n-Propylamii Hexachloroethane	ie					1,2,4,5-Tetrachlorobenzene	
380 U	UG/KG		1			380 U	UG/KG	4-Bromophenyl Phenyl Ether	
380 U	UG/KG	Nitrobenzene				380 U	UG/KG	Hexachlorobenzene (HCB)	· .
. 380 U	UG/KG	Isophorone				380 U	UG/KG	Atrazine	
380 U	UG/KG	2-Nitrophenol				950 U	UG/KG	Pentachlorophenol	
380 U	UG/KG	2,4-Dimethylphenol				380 U	UG/KG	Phenanthrene	•
380 U	UG/KG	bis(2-Chloroethoxy)Meth	ane			380 Ü	UG/KG.	Anthracene	
380 U	UG/KG	2,4-Dichlorophenol				380 U	UG/KG	Carbazole	
380 U	UG/KG	Naphthalene	i			380 U	UG/KG	Di-n-Butylphthalate	
380 U	UG/KG	4-Chloroaniline	ļ.			380 U	UG/KG	Fluoranthene	
380 U	UG/KG	Hexachlorobutadiene				380 U	UG/KG	Pyrene	
380 U	UG/KG	Caprolactam				380 U	UG/KG	Benzyl Butyl Phthalate	
380 U	UG/KG	4-Chloro-3-Methylpheno	1			380 U	UG/KG	3,3'-Dichlorobenzidine	
380 U	UG/KG	2-Methylnaphthalene	(1/000)			380 U	UG/KG	Benzo(a)Anthracene	
380 U	UG/KG	Hexachlorocyclopentadie	ne (HCCP)			380 U	UG/KG	Chrysene	
380 U	UG/KG	2,4,6-Trichlorophenol				380 U	UG/KG	bis(2-Ethylhexyl) Phthalate	
950 U	UG/KG	2,4,5-Trichlorophenol				380 U	UG/KG	Di-n-Octylphthalate	
380 U	UG/KG	1,1-Biphenyl	1		•	380 U	UG/KG	Benzo(b)Fluoranthene	
380 U	UG/KG	2-Chloronaphthalene				380 U	UG/KG	Benzo(k)Fluoranthene	
950 U	UG/KG	2-Nitroaniline	i I			380 U	UG/KG	Benzo-a-Pyrene	
380 U	UG/KG	Dimethyl Phthalate				380 U	UG/KG	Indeno (1,2,3-cd) Pyrene	
380 U	UG/KG	2,6-Dinitrotoluene				380 U	UG/KG	Dibenzo(a,h)Anthracene	
380 U	UG/KG	Acenaphthylene				380 U	UG/KG	Benzo(ghi)Perylene	•
950 U	UG/KG	3-Nitroaniline				13	%	% Moisture	
380 U	UG/KG	Acenaphthene	i						
950 UJ	UG/KG	2,4-Dinitrophenol	!					·	
950 U	UG/KG <sup>-</sup>	4-Nitrophenol	!						
			1						

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | ÚJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is an estimate. | N-Presumptive evidence analyte is present; analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

Project: 03-0474

6794 FY 2003

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Program: SF

Flowood, MS

Id/Station: GS22SS /

MD No: 1XX4

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XX4

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 16:42

Produced by: Goddard, Denise

Ending:

**RESULTS UNITS ANALYTE** 

110 NJ UG/KG HEXADECANOIC ACID

1300 J UG/KG 4 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Ident|fication of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

Sample Project: 03-0474 FY 2003 6795

**Extractables Scan** 

Facility: Gulf States Creosoting

- Flowood, MS

Program: SF

Id/Station: GS06SS /

MD No: 1XX5

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XX5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Endina:

Requestor:

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 16:40

RESULTS	UNITS	ANALYTE		RESULTS	UNITS	ANALYTE
450 U	UG/KG	Benzaldehyde		450 U	UG/KG	Dibenzofuran
450 U	UG/KG	Phenol	,	450 U	UG/KG	2,4-Dinitrotoluene
450 U	UG/KG	bis(2-Chloroethyl) Ether		450 U	UG/KG	Diethyl Phthalate
450 U	UG/KG	2-Chlorophenol		450 U	UG/KG	Fluorene
450 U	· UG/KG	2-Methylphenol		450 U	UG/KG	4-Chlorophenyl Phenyl Ether
450 U	UG/KG	bis(2-Chloroisopropyl) E	ther	1100 U	UG/KG	4-Nitroaniline
450 U	UG/KG	Acetophenone		1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
450 U	UG/KG	(3-and/or 4-)Methylphen		450 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
450 U	UG/KG	n-Nitrosodi-n-Propylamir	ije	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
450 U	UG/KG	Hexachloroethane	! ;	450 U	UG/KG	4-Bromophenyl Phenyl Ether
450 U	UG/KG	Nitrobenzene	Į Į	450 Û	UG/KG	Hexachlorobenzene (HCB)
450 U	UG/KG	Isophorone		450 U	UG/KG	Atrazine
450 U	UG/KG	2-Nitrophenol		1100 U	UG/KG	Pentachlorophenol
450 U 🕟	UG/KG	2,4-Dimethylphenol	!	450 U	UG/KG	Phenanthrene
450 U	UG/KG	bis(2-Chloroethoxy)Meth	ane	450 U	UG/KG	Anthracene
450 U	UG/KG	2,4-Dichlorophenol		450 U	UG/KG	Carbazole
450 U	UG/KG	Naphthalene		450 U	UG/KG	Di-n-Butylphthalate
450 U	UG/KG	4-Chloroaniline		450 U	UG/KG	Fluoranthene
450 U	UG/KG	Hexachlorobutadiene	ļ	450 U	UG/KG	Pyrene
450 U	UG/KG	Caprolactam	·	450 U	UG/KG	Benzyl Butyl Phthalate
450 U	UG/KG	4-Chloro-3-Methylphenol		450 U	UG/KG	3,3'-Dichlorobenzidine
450 U	UG/KG	2-Methylnaphthalene	(11000)	450 U	UG/KG	Benzo(a)Anthracene
450 UJ	UG/KG	Hexachlorocyclopentadie	ene (HCCP)	450 U	UG/KG	Chrysene
450 U	UG/KG	2,4,6-Trichlorophenol	·	450 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol		450 U	UG/KG	Di-n-Octylphthalate
450 U	UG/KG	1,1-Biphenyl		450 U	UG/KG	Benzo(b)Fluoranthene
450 U	UG/KG	2-Chloronaphthalene 2-Nitroaniline		450 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG			450 U	UG/KG	Benzo-a-Pyrene
450 U	UG/KG	Dimethyl Phthalate 2.6-Dinitrotoluene		450 U	UG/KG	Indeno (1,2,3-cd) Pyrene
450 U	UG/KG			450 U	UG/KG	Dibenzo(a,h)Anthracene
450 U	UG/KG	Acenaphthylene 3-Nitroaniline		450 U	UG/KG	Benzo(ghi)Perylene
	UG/KG	=		26	%	% Moisture
	UG/KG	Acenaphthene		•		
	UG/KG	2,4-Dinitrophenol 4-Nitrophenol				
1100 0	UG/KG	4-Millophenoi				
		•				•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6795

FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS06SS /

MD No: 1XX5

Media: SURFACE SOIL (0" - 12")

Case No: 31635

D No: 1XX5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:40

Ending:

**ANALYTE RESULTS UNITS** 100 NJ

UG/KG 170 NJ UG/KG UG/KG HEXADECANOIC ACID 9-OCTADECENE, (E) -1-HEPTADECENE

120 NJ UG/KG 860 NJ 460 NJ UG/KG 2800 J UG/KG

D:C-FRIEDOOLEANAN-3-ONE STIGMAST-4-EN-3-ONE 14 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

450 U

450 U

1100 U

450 U

1100 UJ

1100 U

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

2,6-Dinitrotoluene

Acenaphthylene

3-Nitroaniline

4-Nitrophenol

Acenaphthene

2.4-Dinitrophenol

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise 6796 FY 2003 Project: 03-0474 Sample Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 16:50 Program: SF Case No: 31635 Ending: Id/Station: GS06SB / MD No: 1XX6 Inorg Contractor: SENTIN Org Contractor: LIBRTY D No: 1XX6 Media: SUBSURFACE SOIL (> 12") DATA REPORTED ON DRY WEIGHT BASIS **ANALYTE RESULTS UNITS ANALYTE RESULTS UNITS** 450 U UG/KG 450 U UG/KG Benzaldehyde Dibenzofuran 450 U UG/KG Phenol 450 U UG/KG 2,4-Dinitrotoluene bis(2-Chloroethyl) Ether 450 U UG/KG Diethyl Phthalate 450 U UG/KG 2-Chlorophenol 450 U UG/KG Fluorene 450 U UG/KG 450 U 2-Methylphenol 450 Ù UG/KG 4-Chlorophenyl Phenyl Ether UG/KG bis(2-Chloroisopropyl) Ether 1100 U UG/KG 4-Nitroaniline 450 U UG/KG 1100 UJ 2-Methyl-4,6-Dinitrophenol 450 U UG/KG Acetophenone UG/KG n-Nitrosodiphenylamine/Diphenylamine (3-and/or 4-)Methylphenol 450 U UG/KG 450 U UG/KG UG/KG n-Nitrosodi-n-Propylamine NA UG/KG 1,2,4,5-Tetrachlorobenzene 450 U 450 U UG/KG Hexachloroethane 450 U UG/KG 4-Bromophenyl Phenyl Ether Nitrobenzene 450 U UG/KG Hexachlorobenzene (HCB) 450 U UG/KG 450 U UG/KG Isophorone 450 U UG/KG Atrazine 450 U 2-Nitrophenol 1100 U UG/KG Pentachlorophenol UG/KG UG/KG 450 U UG/KG 450 U 2,4-Dimethylphenol Phenanthrene 450 U UG/KG bis(2-Chloroethoxy)Methane 450 U UG/KG Anthracene 450 U UG/KG 2.4-Dichlorophenol 450 U UG/KG Carbazole 450 U UG/KG Naphthalene 450 U UG/KG Di-n-Butylphthalate 4-Chloroaniline 450 U 450 U UG/KG UG/KG Fluoranthene 450 U UG/KG Hexachlorobutadiene 450 U UG/KG Pyrene 450 U - UG/KG Caprolactam 450 U UG/KG Benzyl Butyl Phthalate 450 U 4-Chloro-3-Methylphenol UG/KG 450 U UG/KG 3,3'-Dichlorobenzidine UG/KG 450 U 2-Methylnaphthalene 450 U UG/KG Benzo(a)Anthracene Hexachlorocyclopentadiene (HCCP) 450 U UG/KG 450 U UG/KG Chrysene 450 U UG/KG 2,4,6-Trichlorophenol 450 U UG/KG bis(2-Ethylhexyl) Phthalate 1100.U UG/KG 2,4,5-Trichlorophenol 450 U UG/KG Di-n-Octylphthalate 450 U UG/KG 1,1-Biphenyl 450 U UG/KG Benzo(b)Fluoranthene 450 U UG/KG 2-Chloronaphthalene 450 U Benzo(k)Fluoranthene UG/KG 1100 U UG/KG 2-Nitroaniline 450 U UG/KG Benzo-a-Pyrene UG/KG Dimethyl Phthalate 450 U 450 U UG/KG Indeno (1,2,3-cd) Pyrene

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

450 U

450 U

26

UG/KG

UG/KG

%

Dibenzo(a,h)Anthracene

Benzo(ghi)Pervlene

% Moisture

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

6797 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:35 Program: SF Case No: 31635 Endina: ld/Station: GS24SS / Inorg Contractor: SENTIN MD No: 1XX7 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX7 DATA REPORTED ON DRY WEIGHT BASIS RESULTS UNITS **ANALYTE** RESULTS UNITS **ANALYTE** 600 U UG/KG 600 U UG/KG Benzaldehyde Dibenzofuran 600 U UG/KG 2.4-Dinitrotoluene 600 U UG/KG Phenol bis(2-Chloroethyl) Ether 600 U Diethyl Phthalate 600 U UG/KG UG/KG 600 U UG/KG 2-Chlorophenol 600 U UG/KG Fluorene 600 U . UG/KG 2-Methylphenol 600 U UG/KG 4-Chlorophenyl Phenyl Ether bis(2-Chloroisopropyl) Ether 1500 U 600 U UG/KG UG/KG 4-Nitroaniline 1500 UJ UG/KG Acetophenone UG/KG 2-Methyl-4,6-Dinitrophenol 600 U 600 U UG/KG (3-and/or 4-)Methylphenol 600 U UG/KG n-Nitrosodiphenylamine/Diphenylamine 600 U UG/KG n-Nitrosodi-n-Propylamine NA UG/KG 1.2.4.5-Tetrachlorobenzene 600 U UG/KG Hexachloroethane 600 U UG/KG 4-Bromophenyl Phenyl Ether 600 U UG/KG 600 U Hexachlorobenzene (HCB) Nitrobenzene UG/KG 600 U UG/KG Isophorone 600 U UG/KG Atrazine 600 U UG/KG 2-Nitrophenol 1500 U UG/KG Pentachlorophenol 600 U UG/KG 2,4-Dimethylphenol 600 U UG/KG Phenanthrene 600 U UG/KG bis(2-Chloroethoxy)Methane 330 J UG/KG Anthracene 600 U UG/KG 2.4-Dichlorophenol 600 U UG/KG Carbazole 600 U UG/KG Naphthalene 600 U UG/KG Di-n-Butylphthalate 600 U ·UG/KG 4-Chloroaniline 1300 UG/KG Fluoranthene UG/KG 600 U Hexachlorobutadiene UG/KG Pyrene 1600 600 U UG/KG Caprolactam 600 U UG/KG Benzyl Butyl Phthalate 600 U UG/KG 4-Chloro-3-Methylphenol 600 U UG/KG 3.3'-Dichlorobenzidine 600 U UG/KG 2-Methylnaphthalene 990 UG/KG Benzo(a)Anthracene 600 U UG/KG Hexachlorocyclopentadiene (HCCP) 1100 UG/KG Chrysene 600 U UG/KG 2,4,6-Trichlorophenol 600 U UG/KG bis(2-Ethylhexyl) Phthalate 1500 U UG/KG 2,4,5-Trichlorophenol 600 U UG/KG Di-n-Octvlphthalate 600 U UG/KG 1.1-Biphenyl 2700 J UG/KG Benzo(b)Fluoranthene 2-Chloronaphthalene 600 U UG/KG 2600 J UG/KG Benzo(k)Fluoranthene 1500 U UG/KG 2-Nitroaniline 870 UG/KG Benzo-a-Pyrene 600 U UG/KG Dimethyl Phthalate 550 J UG/KG Indeno (1,2,3-cd) Pyrene 600 U UG/KG 2.6-Dinitrotoluene 260 J UG/KG Dibenzo(a,h)Anthracene UG/KG 150 J Acenaphthylene 330 J UG/KG Benzo(ghi)Perylene 1500 U UG/KG 3-Nitroaniline 45 % % Moisture 600 U TUG/KG Acenaphthene 1500 UJ UG/KG 2,4-Dinitrophenol 1500 U UG/KG 4-Nitrophenol

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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# **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6797 FY 2003

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Program: SF ld/Station: GS24SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635 MD No: 1XX7

D No: 1XX7

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:35

Ending:

**RESULTS UNITS ANALYTE** 1600 NJ UG/KG **CARYOPHYLLENE** UG/KG 160 NJ PYRENE, 1-METHYL-7H-BENZ [DE] ANTHRACEN-7-ONE 140 NJ UG/KG BENZO [E] PYRENE 840 NJ UG/KG **UNKNOWN ALKENE** 200 J UG/KG 2600 NJ UG/KG STIGMAST-4-EN-3-ONE UG/KG 1700 J **6 UNKNOWN PAHS** 5900 J UG/KG 15 UNKNOWN COMPOUNDS 260 UG/KG **PERYLENE** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6798 FY 2003

Produced by: Goddard, Denise

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO

**Extractables Scan** 

Case No: 31635

Beginning: 04/23/2003 07:40

Program: SF

Ending:

Id/Station: GS12SS /

MD No: 1XX9

**D No: 1XX9** 

Flowood, MS

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
400 U	UG/KG	Benzaldehyde	400 U	UG/KG	Dibenzofuran
400 U	UG/KG	Phenol	400 U	UG/KG .	2,4-Dinitrotoluene
400 U	UG/KG	bis(2-Chloroethyl) Ether	400 U	UG/KG	Diethyl Phthalate
400 U	UG/KG	2-Chlorophenol	400 U	UG/KG	Fluorene
400 U	UG/KG	2-Methylphenol	400 U	UG/KG	4-Chlorophenyl Phenyl Ether
400 U	UG/KG	bis(2-Chloroisopropyl) Ether	1000 U	UG/KG	4-Nitroaniline
400 U	UG/KG	Acetophenone	1000 UJ	UG/KG	2-Methyl-4,6-Dinitrophenol
400 U	UG/KG	(3-and/or 4-)Methylphenol	400 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
400 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
400 U	UG/KG	Hexachloroethane	400 U	UG/KG	4-Bromophenyl Phenyl Ether
400 U	UG/KG	Nitrobenzene	400 U	UG/KG	Hexachlorobenzene (HCB)
400 U	UG/KG	Isophorone	400 U	UG/KG	Atrazine
400 U	UG/KG	2-Nitrophenol	1000 U	UG/KG	Pentachlorophenol
400 U	UG/KG	2,4-Dimethylphenol	400 U	UG/KG	Phenanthrene
	* UG/KG	bis(2-Chloroethoxy)Methane	400 U	UG/KG	Anthracene
400 U	UG/KG	2,4-Dichlorophenol	400 U	UG/KG	Carbazole
400 U	UG/KG	Naphthalene	. 400 U	UG/KG	Di-n-Butylphthalate
400 U	UG/KG	4-Chloroaniline	400 U	UG/KG	Fluoranthene
400 U	UG/KG	Hexachlorobutadiene	400 U	UG/KG	Pyrene
400 U	UG/KG	Caprolactam	400 U	.UG/KG	Benzyl Butyl Phthalate
. 400 U	UG/KG	4-Chloro-3-Methylphenol	400 U	UG/KG	3,3'-Dichlorobenzidine
400 U	UG/KG	2-Methylnaphthalene	400 U	UG/KG	Benzo(a)Anthracene
400 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	400 U	ÚG/KG	Chrysene
400 U	UG/KG	2,4,6-Trichlorophenol	400 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1000 U	UG/KG	2,4,5-Trichlorophenol	400 U	UG/KG	Di-n-Octylphthalate
400 U	UG/KG	1,1-Biphenyl	400 U	UG/KG	Benzo(b)Fluoranthene
400 U	UG/KG	2-Chioronaphthalene	400 U	UG/KG	Benzo(k)Fluoranthene
1000 U	UG/KG	2-Nitroaniline	400 U	UG/KG	Benzo-a-Pyrene
400 U	UG/KG	Dimethyl Phthalate	400 U	UG/KG	Indeno (1,2,3-cd) Pyrene
400 U	UG/KG	2,6-Dinitrotoluene	400 U	UG/KG	Dibenzo(a,h)Anthracene
400 U	UG/KG	Acenaphthylene	400 U	UG/KG	Benzo(ghi)Perylene
1000 U	UG/KG	3-Nitroaniline	17	%	% Moisture
400 U	UG/KG	Acenaphthene			
1000 UJ	UG/KG	2,4-Dinitrophenol			
1000 U	UG/KG	4-Nitrophenol			
					· ·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6798 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO

Beginning: 04/23/2003 07:40

Program: SF

410 J

Case No: 31635

Ending:

Id/Station: GS12SS / Media: SURFACE SOIL (0" - 12") MD No: 1XX9 D No: 1XX9

Flowood, MS

Inorg Contractor: SENTIN Org Contractor: LIBRTY

RESULTS UNITS

**ANALYTE** 

UG/KG

3 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:10

6799 FY 2003 Project: 03-0474 Sample

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS12SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XY0 D No: 1XY0

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/23/2003 07:50

Requestor:

Ending:

<b>RESULTS</b>	UNITS	ANALYTE	: ••	RESULTS	UNITS	ANALYTE
450 U	UG/KG	Benzaldehyde	;	450 U	UG/KG	Dibenzofuran
450 U	UG/KG	Phenol		450 U	UG/KG	2,4-Dinitrotoluene
450 U	UG/KG	bis(2-Chloroethyl) Ether	· <b>f</b>	450 U	UG/KG	Diethyl Phthalate
450 U	UG/KG	2-Chlorophenol		450 U	UG/KG	Fluorene
450 U	UG/KG	2-Methylphenol		450 U	UG/KG	4-Chlorophenyl Phenyl Ether
450 U	UG/KG	bis(2-Chloroisopropyl) E	ther	1100 U	UG/KG	4-Nitroaniline
450 U	UG/KG	Acetophenone	·	1100 UJ	UG/KG	2-Methyl-4,6-Dinitrophenol
450 U	UG/KG	(3-and/or 4-)Methylpher		450 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
450 U	UG/KG	n-Nitrosodi-n-Propylami	ine .		UG/KG	1,2,4,5-Tetrachlorobenzene
450 U	UG/KG	Hexachloroethane	·	450 U	UG/KG	4-Bromophenyl Phenyl Ether
450 U	UG/KG	Nitrobenzene		450 U	UG/KG	Hexachlorobenzene (HCB).
450 U	UG/KG	Isophorone	'	450 U	UG/KG	Atrazine
450 U	UG/KG	2-Nitrophenol		1100 U	UG/KG	Pentachlorophenol
450 U	UG/KG	2,4-Dimethylphenol		450 U	UG/KG	Phenanthrene
450 U	UG/KG	bis(2-Chloroethoxy)Met	hane	450 U	UG/KG	Anthracene
450. U	UG/KG	2,4-Dichlorophenol		450 U	UG/KG	Carbazole
450 U	UG/KG	Naphthalene		450 U	UG/KG	Di-n-Butylphthalate
450 U	UG/KG	4-Chloroaniline		450 U	UG/KG	Fluoranthene
_ 450 U	UG/KG	Hexachlorobutadiene		450 <sub>.</sub> U	UG/KG	Pyrene
450 U	UG/KG	Caprolactam		450 U	UG/KG	Benzyl Butyl Phthalate
450 U	UG/KG	4-Chloro-3-Methylpheno	اَمْ	450 U	UG/KG	3,3'-Dichlorobenzidine
450 U	UG/KG	2-Methylnaphthalene		450 U	UG/KG	Benzo(a)Anthracene
450 U	UG/KG	Hexachlorocyclopentadi	iene (HCCP)	450 U	UG/KG	Chrysene
	UG/KG	2,4,6-Trichlorophenol		450 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol		450 U	UG/KG	Di-n-Octylphthalate
450 U	UG/KG	1,1-Biphenyl		450 U	UG/KG	Benzo(b)Fluoranthene
450 U	UG/KG	2-Chloronaphthalene		450 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline		450 U	UG/KG	Benzo-a-Pyrene
450 U	UG/KG	Dimethyl Phthalate		450 U	UG/KG	Indeno (1,2,3-cd) Pyrene
450 U	UG/KG	2,6-Dinitrotoluene		450 U	UG/KG	Dibenzo(a,h)Anthracene
450 U	UG/KG	Acenaphthylene		450 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline		26	%	% Moisture
450 U	UG/KG	Acenaphthene			•	
1100 UJ	UG/KG	2,4-Dinitrophenol				
1100 U	UG/KG	4-Nitrophenot				
		İ				
			İ	•		·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise Sample 6800 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 07:40 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS03SS / MD No: 1XY1 Org Contractor: LIBRTY. Media: SURFACE SOIL (0" - 12") D No: 1XY1 DATA REPORTED ON DRY WEIGHT BASIS

			<u>:</u>						·		
RESULTS	UNITS	ANALYTE	: !. !					RESULTS	UNITS	ANALYTE	
430 U	UG/KG	Benzaldehyde	! :					430 U	UG/KG	Dibenzofuran	
430 U	UG/KG	Phenol	İ					430 U	UG/KG	2,4-Dinitrotoluene	
430 U	UG/KG	bis(2-Chloroethyl) Ether						430 U	UG/KG	Diethyl Phthalate	
430 U	UG/KG	2-Chlorophenol						430 U	UG/KG	Fluorene	
430 U	UG/KG	2-Methylphenol	:		•			430 U	UG/KG	4-Chlorophenyl Phenyl Ether	
430 U	UG/KG	bis(2-Chloroisopropyl) E	ther					1100 U	UG/KG	4-Nitroaniline	
430 U	UG/KG	Acetophenone						1100 UJ	UG/KG	2-Methyl-4,6-Dinitrophenol	i
430 U	UG/KG	(3-and/or 4-)Methylpher	iol					430 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine	
. 430 U	UG/KG	n-Nitrosodi-n-Propylami	ne			•		NA	UG/KG	1,2,4,5-Tetrachlorobenzene	
430 U	UG/KG	Hexachloroethane	1					430 U	UG/KG	4-Bromophenyl Phenyl Ether	
430 U	UG/KG	Nitrobenzene	!					430 U	UG/KG	Hexachlorobenzene (HCB)	
430 U	UG/KG	Isophorone						430 U	UG/KG	Atrazine	
430 U	UG/KG	2-Nitrophenol	į	•				1100 U	.UG/KG	Pentachlorophenol	
430 U	UG/KG	2,4-Dimethylphenol	!					430 U	UG/KG 1	Phenanthrene	
430 U	UG/KG	bis(2-Chloroethoxy)Metl	hane					430 U	UG/KG	Anthracene.	
430 U	UG/KG	2,4-Dichlorophenol						430 U	UG/KG	Carbazole	
430 U	UG/KG	Naphthalene	!		•			430 U	UG/KG	Di-n-Butylphthalate	
430 U	UG/KG	4-Chloroaniline	į					430 U	'UG/KG	Fluoranthene	
430 U	UG/KG	Hexachlorobutadiene				•		430 U	UG/KG	Pyrene	
430 U	UG/KG	Caprolactam						430 U	UG/KG	Benzyl Butyl Phthalate	
430 U	UG/KG	4-Chloro-3-Methylpheno	il					430 U	UG/KG	3,3'-Dichlorobenzidine	
430 U	UG/KG	2-Methylnaphthalene	į.					430 U	UG/KG	Benzo(a)Anthracene	
430 U	UG/KG	Hexachlorocyclopentadi	ene (HCCP)			•	٠.	430 U	UG/KG	Chrysene	
430 U	UG/KG	2,4,6-Trichlorophenol	!					430 U	UG/KG	bis(2-Ethylhexyl) Phthalate	
1100 U	UG/KG	2,4,5-Trichlorophenol						430 U	UG/KG	Di-n-Octylphthalate	
-430 U	UG/KG	1,1-Biphenyl	; !					420 U	UG/KG	Benzo(b)Fluoranthene	•
430 U	UG/KG	2-Chloronaphthalene	!					430 U	UG/KG	Benzo(k)Fluoranthene	
1100 U	UG/KG	2-Nitroaniline						430 U	UG/KG	Benzo-a-Pyrene	
430 U	UG/KG	Dimethyl Phthalate						430 U	UG/KG	Indeno (1,2,3-cd) Pyrene	•
430 U	UG/KG	2,6-Dinitrotoluene		•				430 U	UG/KG	Dibenzo(a,h)Anthracene	
430 U	UG/KG	Acenaphthylene	!		'			430 U	UG/KG	Benzo(ghi)Perylene	•
1100 U	UG/KG	3-Nitroaniline						24	%	% Moisture	
430 U	UG/KG	Acenaphthene	1								
1100 UJ	UG/KG	2,4-Dinitrophenol									
1100 U	UG/KG	4-Nitrophenol	2							•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD. ATHENS. GA** 

Production Date: 06/20/2003 14:10

Sample

6800 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 07:40

Ending:

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS03SS /

MD No: 1XY1

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

D No: 1XY1

Org Contractor: LIBRTY

**RESULTS UNITS** 

**ANALYTE** 

91 J UG/KG **UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6801 FY 2003

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS03SB /

MD No: 1XY2

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

D No: 1XY2

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Ending:

Requestor:

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/23/2003 07:55

			<del></del>	 				
RESULTS	UNITS	ANALYTE				RESULTS	UNITS	ANALYTE
460 U	UG/KG	Benzaldehyde				460 U	UG/KG	Dibenzofuran
460 U	UG/KG	Phenol	1			460 U	UG/KG	2,4-Dinitrotoluene
460 U	UG/KG	bis(2-Chloroethyl) Ether				460 U	UG/KG	Diethyl Phthalate
460 U	UĠ/KG	2-Chlorophenol	<i>i</i> :			460 U	UG/KG	Fluorene
460 U	UG/KG	2-Methylphenol	i		•	460 U	UG/KG	4-Chlorophenyl Phenyl Ether
460 U	UG/KG	bis(2-Chloroisopropyl) E	ther			1200 U	UG/KG	4-Nitroaniline
460 U	UG/KG	Acetophenone				1200 U	UG/KG	2-Methyl-4,6-Dinitrophenol
460 U	UG/KG	(3-and/or 4-)Methylpher	ol			460 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
460 U	UG/KG	n-Nitrosodi-n-Propylami	ne			NA	UG/KG	1,2,4,5-Tetrachlorobenzene
460 U	UG/KG	Hexachloroethane				460 U	ŪG/KG	4-Bromophenyl Phenyl Ether
460 U	UG/KG	Nitrobenzene				460 U	UG/KG	Hexachlorobenzene (HCB)
460 U	UG/KG	Isophorone				460 U	UG/KG	Atrazine
460 U	UG/KG	2-Nitrophenol	į			1200 U	UG/KG	Pentachlorophenol
460 U	UG/KG	2,4-Dimethylphenol	!			460 U	UG/KG	Phenanthrene
460 U	UG/KG	bis(2-Chloroethoxy)Meth	iane			460 U	UG/KG	Anthracene
460 U	UG/KG	2,4-Dichlorophenol	!			460 U	UG/KG	Carbazole
460 U	UG/KG	Naphthalene			•	460 U	UG/KG	Di-n-Butylphthalate
460 U	UG/KG	4-Chloroaniline				460 U .	UG/KG	Fluoranthene
460 U	UG/KG	Hexachlorobutadiene	i I			460 U	UG/KG	Pyrene
460 U	UG/KG	Caprolactam	! !			460 UJ	UG/KG	Benzyl Butyl Phthalate
460 U	UG/KG	4-Chloro-3-Methylpheno	1			460 U	UG/KG	3,3'-Dichlorobenzidine
460 U	UG/KG	2-Methylnaphthalene				460 U	UG/KG	Benzo(a)Anthracene
460 U	UG/KG	Hexachlorocyclopentadi	ene (HCCP)			460 U	UG/KG	Chrysene
460 U	UG/KG	2,4,6-Trichlorophenol				460 UJ	UG/KG	bis(2-Ethylhexyl) Phthalate
1200 U .	UG/KG	2,4,5-Trichlorophenol	;			460 U	UG/KG	Di-n-Octylphthalate
460 U	UG/KG	1,1-Biphenyl				460 U	UG/KG	Benzo(b)Fluoranthene
460 U	UG/KG	2-Chloronaphthalene	i			460 U	UG/KG	Benzo(k)Fluoranthene
1200 U	UG/KG	2-Nitroaniline				460 U	UG/KG	Benzo-a-Pyrene
460 U	UG/KG	Dimethyl Phthalate				460 U	UG/KG	Indeno (1,2,3-cd) Pyrene
460 U	UG/KG	2,6-Dinitrotoluene	i	•		460 U	UG/KG	Dibenzo(a,h)Anthracene
460 U-	UG/KG	Acenaphthylene				460 U	UG/KG	Benzo(ghi)Perylene
1200 U	UG/KG	3-Nitroaniline	į.			28	%	% Moisture
460 U	UG/KG	Acenaphthene				•		
1200 U	UG/KG	2,4-Dinitrophenol						
1200 U	UG/KG	4-Nitrophenol			•			
		•						

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**MISCELLANEOUS COMPOUNDS** 

Media: SUBSURFACE SOIL (> 12")

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/23/2003 07:55

6801

Requestor:

Program: SF

Facility: Gulf States Creosoting

Flowood, MS

Id/Station: GS03SB /

Case No: 31635 MD No: 1XY2

D No: 1XY2

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Ending:

**RESULTS UNITS** 

ANALYTE

250 J UG/KG 2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6802 FY 2003 Project: 03-0474

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS08SS /

**Extractables Scan** 

Media: SURFACE SOIL (0" - 12")

Case No: 31635

MD No: 1XY3 D No: 1XY3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:10

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE
1400 U	UG/KG	Benzaldehyde			1400 U	UG/KG	Dibenzofuran
1400 U	UG/KG	Phenol	•		1400 U	UG/KG	2,4-Dinitrotoluene
1400 U	UG/KG	bis(2-Chloroethyl) Ether			1400 U	UG/KG	Diethyl Phthalate
1400 U	UG/KG	2-Chlorophenol	• •		1400 U	UG/KG	Fluorene
1400 U	UG/KG	2-Methylphenol			1400 U	UG/KG	4-Chlorophenyl Phenyl Ether
1400 U	UG/KG	bis(2-Chloroisopropyl) Ether			3500 U	UG/KG	4-Nitroaniline
1400 U -	UG/KG	Acetophenone	_		3500 UJ	UG/KG	2-Methyl-4,6-Dinitrophenol
1400 U	UG/KG	(3-and/or 4-)Methylphenol			1400 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
1400 U	UG/KG `	n-Nitrosodi-n-Propylamine			NA	UG/KG	1,2,4,5-Tetrachlorobenzene
1400 U	UG/KG	Hexachloroethane			1400 U	UG/KG	4-Bromophenyl Phenyl Ether
1400 U	UG/KG	Nitrobenzene			1400 U	UG/KG	Hexachlorobenzene (HCB)
. 1400 U	UG/KG	Isophorone	•		1400 U	UG/KG	Atrazine
1400 U	UG/KG	2-Nitrophenol	•		3500 U	UG/KG	Pentachlorophenol
1400 U	UG/KG	2,4-Dimethylphenol	•		-610 J	UG/KG	Phenanthrene
1400 U	UG/KG	bis(2-Chloroethoxy)Methane	•		930 J	UG/KG	Anthracene
1400 U	UG/KG	2,4-Dichlorophenol			190 J	UG/KG	Carbazole
1400 U	UG/KG	Naphthalene			1400 U	UG/KG	Di-n-Butylphthalate
1400 U	UG/KG	4-Chloroaniline		•	2600	UG/KG	Fluoranthene
1400 U	UG/KG	Hexachlorobutadiene			2500	UG/KG	Pyrene
1400 U	UG/KG	Caprolactam		,	1400 U	UG/KG	Benzyl Butyl Phthalate
1400 U	UG/KG	4-Chloro-3-Methylphenol	•	•	1400 U	UG/KG	3,3'-Dichlorobenzidine
1400 U	UG/KG	2-Methylnaphthalene			1500	UG/KG	Benzo(a)Anthracene
1400 U	UG/KG	Hexachlorocyclopentadiene (HCCP)			2300	UG/KG	Chrysene
1400 U	UG/KG	2,4,6-Trichlorophenol			1400 U	UG/KG	bis(2-Ethylhexyl) Phthalate
3500 U	UG/KG	2,4,5-Trichlorophenol			1400 U	UG/KG	Di-n-Octylphthalate
1400 U	UG/KG	1,1-Biphenyl			11000 J	UG/KG	Benzo(b)Fluoranthene
1400 U	UG/KG	2-Chloronaphthalene			12000 J	UG/KG	Benzo(k)Fluoranthene
3500 U	UG/KG	2-Nitroaniline			3700	UG/KG	Benzo-a-Pyrene
1400 U	UG/KG	Dimethyl Phthalate		٠.	5000	UG/KG	Indeno (1,2,3-cd) Pyrene
1400 U	UG/KG	2,6-Dinitrotoluene			1700	UG/KG	Dibenzo(a,h)Anthracene
480 J	UG/KG	Acenaphthylene			3800	UG/KG	Benzo(ghi)Perylene
3500 U	UG/KG	3-Nitroaniline	•		4	%	% Moisture
1400 U	UG/KG .	Acenaphthene	•				·
3500 UJ	UG/KG	2,4-Dinitrophenol					
3500 U *	UG/KG	4-Nitrophenol					
			•			•	
							•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable, reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable, reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample 6802 FY 2003 Project: 03-0474 Produced by: Goddard, Denise

**MISCELLANEOUS COMPOUNDS** 

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO

Program: SF

Flowood, MS

Beginning: 04/23/2003 08:10

Case No: 31635

Ending:

Id/Station: GS08SS /

MD No: 1XY3

D No: 1XY3

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

Org Contractor: LIBRTY

RESULTS	UNITS	ANALYTE
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	•	
920 J	UG/KG	BENZOFLUORENE
1000 NJ	UG/KG	PYRENE, 2-METHYL-
880_NJ	UG/KG	PYRENE, 1-METHYL-
910 NJ	UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
1200 NJ	UG/KG	CHRYSENE, 1-METHYL-
1400 NJ	UG/KG	5,12-NAPHTHACENEDIONE
3900 NJ	UG/KG	PERYLENE
880 NJ	UG/KG	1,2:4,5-DIBENZPYRENE
5400 J	UG/KG	4 UNKNOWN PAHS
3000 1	LIG/KG	17 LINKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6804 FY **2003** Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 08:20 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS08SB / MD No: 1XY4 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XY4 DATA REPORTED ON DRY WEIGHT BASIS **ANALYTE RESULTS UNITS** ANALYTE RESULTS UNITS 400 U UG/KG 400 U UG/KG Benzaldehyde Dibenzofuran 400 U 400 U UG/KG Phenol UG/KG 2,4-Dinitrotoluene 400 U UG/KG bis(2-Chloroethyl) Ether 400 U UG/KG Diethyl Phthalate 400 U UG/KG 2-Chlorophenol Fluorene 400 U UG/KG 400 U UG/KG 4-Chlorophenyl Phenyl Ether 400 U UG/KG 2-Methylphenol 400 U UG/KG bis(2-Chloroisopropyl) Ether 1000 U UG/KG 4-Nitroaniline 400 U 1000 UJ UG/KG 2-Methyl-4,6-Dinitrophenol UG/KG Acetophenone 400 U 400 U UG/KG (3-and/or 4-)Methylphenol UG/KG n-Nitrosodiphenylamine/Diphenylamine n-Nitrosodi-n-Propylamine UG/KG 1.2.4.5-Tetrachlorobenzene 400 U UG/KG NA 400 U UG/KG 4-Bromophenyl Phenyl Ether 400 U UG/KG Hexachloroethane 400 U UG/KG Nitrobenzene 400 U UG/KG Hexachlorobenzene (HCB) 400 U UG/KG Isophorone 400 U UG/KG Atrazine 400 U 1000 U UG/KG 2-Nitrophenol UG/KG Pentachlorophenol 400 U UG/KG 2.4-Dimethylphenol 400 U UG/KG Phenanthrene 400 U UG/KG bis(2-Chloroethoxy)Methane 74 J UG/KG Anthracene 400 U. 2.4-Dichlorophenol UG/KG 400 U UG/KG Carbazole 400 U Naphthalene 400 U UG/KG UG/KG Di-n-Butylphthalate 400 U UG/KG 4-Chloroaniline 200 J UG/KG Fluoranthene 400 U UG/KG Hexachlorobutadiene 230 J UG/KG Pyrene 400 U UG/KG Caprolactam 400 U UG/KG Benzyl Butyl Phthalate 400 U 400 U UG/KG 4-Chloro-3-Methylphenol UG/KG 3,3'-Dichlorobenzidine 400 U UG/KG 2-Methylnaphthalene 120 J UG/KG Benzo(a)Anthracene 400 U UG/KG Hexachlorocyclopentadiene (HCCP) 220 J UG/KG Chrysene 400 U UG/KG 2,4,6-Trichlorophenol 400 U UG/KG bis(2-Ethylhexyl) Phthalate 1000 U UG/KG 2,4,5-Trichlorophenol 400 U UG/KG Di-n-Octylphthalate 400 U UG/KG 1.1-Biphenyl 1000 J UG/KG Benzo(b)Fluoranthene 400 U UG/KG 2-Chloronaphthalene 980 J UG/KG Benzo(k)Fluoranthene 1000 U UG/KG 2-Nitroaniline 240 J UG/KG Benzo-a-Pyrene 400 U UG/KG Dimethyl Phthalate 420 UG/KG Indeno (1,2,3-cd) Pyrene 400 U UG/KG 2,6-Dinitrotoluene 150 J UG/KG Dibenzo(a,h)Anthracene 400 U UG/KG Acenaphthylene 320 J UG/KG Benzo(ghi)Perylene 1000 U UG/KG 3-Nitroaniline 17 % Moisture 400 U UG/KG Acenaphthene 1000 UJ UG/KG 2.4-Dinitrophenol 1000 U UG/KG 4-Nitrophenol

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

MISCELLANEOUS COMPOUNDS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6804 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/23/2003 08:20

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS08SB /

MD No: 1XY4

D No: 1XY4

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

Facility: Gulf States Creosoting

Org Contractor: LIBRTY

Ending:

**RESULTS UNITS ANALYTE** 

110 J UG/KG **UNKNOWN AMIDE** 

490 NJ UG/KG 190 NJ UG/KG PERYLENE 1,2:3,4-DIBENZPYRENE

120 NJ UG/KG

NAPHTHO [1,2,3,4-DEF] CHRYSENE

560 J 1500 J

UG/KG UG/KG

3 UNKNOWN PAHS 7 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample 6805 FY 2003 Project: 03-0474 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/23/2003 08:35 Program: SF Case No: 31635 Ending: Id/Station: GS07SS / Inorg Contractor: SENTIN MD No: 1XY5. Media: SURFACE SOIL (0" - 12") Org Contractor: LIBRTY D No: 1XY5 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE				RESULTS	UNITS	ANALYTE	
390 U	UG/KG	Benzaldehyde				. 240 J	UG/KG	Dibenzofuran	
390 U	UG/KG	Phenol				390 U	UG/KG	2,4-Dinitrotoluene	
390 U	UG/KG	bis(2-Ghloroethyl) Ether				390 U	UG/KG	Diethyl Phthalate	
390 U	UG/KG	2-Chlorophenol	_			180 J	UG/KG	Fluorene	
390 U	UG/KG	2-Methylphenol				390 U	UG/KG	4-Chlorophenyl Phenyl Ether	
390 U	UG/KG	bis(2-Chloroisopropyl) Ether	•			980 U	UG/KG	4-Nitroaniline	
390 U	UG/KG	Acetophenone				980 UJ	UG/KG	2-Methyl-4,6-Dinitrophenol	
390 U	UG/KG	(3-and/or 4-)Methylphenol	•	•		390 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine	
390 U	UG/KG	n-Nitrosodi-n-Propylamine	•			NA	UG/KG	1,2,4,5-Tetrachlorobenzene	
390 U	UG/KG	Hexachloroethane				390 U	UG/KG	4-Bromophenyl Phenyl Ether	
390 U	UG/KG	Nitrobenzene	•	•		390 U	UG/KG	Hexachlorobenzene (HCB)	
390 U	UG/KG	Isophorone				390 U	UG/KG	Atrazine	
390 U	UG/KG	2-Nitrophenol				980 U	UG/KG	Pentachlorophenol	
390 U	UG/KG	2,4-Dimethylphenol				4200	UG/KG	Phenanthrene	
390 U	UG/KG	bis(2-Chloroethoxy)Methane				970	UG/KG	Anthracene	
390 U	UG/KG	2,4-Dichlorophenol				510	UG/KG	Carbazole	
87 J	UG/KG	Naphthalene				. 390 U	. UG/KG	Di-n-Butylphthalate	<b>%</b>
390 U	UG/KG	4-Chloroaniline	•			8200	UG/KG	Fluoranthene	
390 U	UG/KG	Hexachlorobutadiene				6200	UG/KG	Pyrene	•
390 U	UG/KG	Caprolactam				390 U	UG/KG	Benzyl Butyl Phthalate	
390 U	UG/KG	4-Chloro-3-Methylphenol		-		390 U	UG/KG	3,3'-Dichlorobenzidine	,
120 J	UG/KG	2-Methylnaphthalene				2500	UG/KG	Benzo(a)Anthracene	
390 U	UG/KG	Hexachlorocyclopentadiene (HCCP)				4900	UG/KG	Chrysene	
390 U	UG/KG	2,4,6-Trichlorophenol	•			390 U	UG/KG	bis(2-Ethylhexyl) Phthalate	
980 U	UG/KG	2,4,5-Trichlorophenol				390 U	UG/KG	Di-n-Octylphthalate	
45 J 🕝	UG/KG	1,1-Biphenyl				4600 J	UG/KG	Benzo(b)Fluoranthene	
390 ⊍	UG/KG	2-Chloronaphthalene			٠	5200 J	UG/KG	Benzo(k)Fluoranthene	
980 Ú	UG/KG	2-Nitroaniline				2100	UG/KG	Benzo-a-Pyrene	
390 Ϣ	UG/KG	Dimethyl Phthalate				1100	UG/KG	Indeno (1,2,3-cd) Pyrene	•".
390 Ú	UG/KG	2,6-Dinitrotoluene				520	UG/KG	Dibenzo(a,h)Anthracene	
320 J	UG/KG	Acenaphthylene				690	UG/KG	Benzo(ghi)Perylene	•
980 U	UG/KG	3-Nitroaniline				15	%	% Moisture	
170 J	UG/KG	Acenaphthene							•
980 UJ	UG/KG	2,4-Dinitrophenol							
980 U	UG/KG	4-Nitrophenol						•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Project: 03-0474 Sample 6805 FY 2003

Produced by: Goddard, Denise

**MISCELLANEOUS COMPOUNDS** 

Requestor:

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO

Program: SF

Case No: 31635

Beginning: 04/23/2003 08:35

ld/Station: GS07SS /

MD No: 1XY5

Ending:

Media: SURFACE SOIL (0" - 12")

**D No: 1XY5** 

Inorg Contractor: SENTIN Org Contractor: LIBRTY

RESULTS	UNITS	ANALYTE
460 NJ	UG/KG	9H-FLUOREN-9-ONE
870 NJ	UG/KG	PHENANTHRENE, 2-METHYL-
1000 J	UG/KG	ANTHRACENE, 2-METHYL-
760 NJ	UG/KG	NAPHTHALENE, 2-PHENYL-
890 NJ	UG/KG	9,10-ANTHRACENEDIONE
830 NJ	UG/KG	PHENANTHRENE, 4,5-DIMETHYL-
740 NJ	UG/KG	11H-BENZO [B] FLUORENE
1200 NJ	UG/KG	PYRENE 2-METHYL-
740 NJ	UG/KG	PYRENE, 4-METHYL-
850 NJ	UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
670 NJ	UG/KG	BENZO [B] NAPHTHO (2,1-D) THIPHENE
580 NJ	UG/KG	CHRYSENE, 5-MEHTHYL-
540 NJ	UG/KG	BENZ (A) ANTHRACENE-7,12-DIONE
1200 NJ	UG/KG	BENZO [E] PYRENE
5300 J	UG/KG	7 UNKNOWN PAHS
5900 J	UG/KG	8 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6806 Requestor: **Extractables Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 08:45 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS07SB / MD No: 1XY6 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XY6 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
450 U	UG/KG	Benzaldehyde	450 U	UG/KG	Dibenzofuran
· 450 U	UG/KG	Phenol	450 U	UG/KG	2.4-Dinitrotoluene
	UG/KG	· · · · · · · · · · · · · · · · · · ·			— · ·
450 U		bis(2-Chloroethyl) Ether	450 U	UG/KG	Diethyl Phthalate
450 U	UG/KG	2-Chlorophenol	450 U	UG/KG	Fluorene
450 U	UG/KG	2-Methylphenol	450 U	UG/KG	4-Chlorophenyl Phenyl Ether
450 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
450 U	. UG/KG	Acetophenone	1100 UJ	UG/KG	2-Methyl-4,6-Dinitrophenol
450 U _		(3-and/or 4-)Methylphenol	450 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
450 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
450 U	UG/KG	Hexachloroethane	450 U	UG/KG	4-Bromophenyl Phenyl Ether
450 U	UG/KG	Nitrobenzene	450 U	UG/KG	Hexachlorobenzene (HCB)
450 U	UG/KG	Isophorone	450 U	UG/KG	Atrazine
450 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
450 U	UG/KG	2,4-Dimethylphenol	450 U	UG/KG	Phenanthrene
450 U	UG/KG	bis(2-Chloroethoxy)Methane	450 U	UG/KG	Anthracene
450 U	UG/KG	2,4-Dichlorophenol	450 U	UG/KG	Carbazole
450 U	UG/KG	Naphthalene	450 U	UG/KG	Di-n-Butylphthalate
450 U	UG/KG	4-Chloroaniline	450 U	UG/KG	Fluoranthene
450 U	UG/KG	Hexachlorobutadiene	450 U	UG/KG	Pyrene
450 U	UG/KG	Caprolactam	450 U	UG/KG	Benzyl Butyl Phthalate
450 U	UG/KG	4-Chloro-3-Methylphenol	450 U	UG/KG	3.3'-Dichlorobenzidine
450 U	UG/KG	2-Methylnaphthalene	450 U	UG/KG	Benzo(a)Anthracene
450 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	450 U	UG/KG	Chrysene
450 U	UG/KG	2,4,6-Trichlorophenol	450 U	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol	450 U	UG/KG	Di-n-Octylphthalate
450 U	UG/KG	1,1-Biphenyl	47 J	UG/KG	Benzo(b)Fluoranthene
450 U	UG/KG	2-Chloronaphthalene	45 J	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	450 U	UG/KG	Benzo-a-Pyrene
450 U	UG/KG	Dimethyl Phthalate	450 U	UG/KG	Indeno (1,2,3-cd) Pyrene
450 Ú	UG/KG	2,6-Dinitrotoluene	450 U `		Dibenzo(a,h)Anthracene
450 U	UG/KG	Acenaphthylene	450 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	26	%	% Moisture
450 U	UG/KG	Acenaphthene			
1100 UJ	UG/KG	2,4-Dinitrophenol			•
1100 U	UG/KG	4-Nitrophenol	•		
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U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be ress than the reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6806

FY **2003** 

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS07SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635 MD No: 1XY6

D No: 1XY6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:45

Ending:

**RESULTS UNITS** 

**ANALYTE** 

190 J UG/KG **UNKNOWN COMPOUND** 

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Inorg Contractor: SENTIN

Id/Station: GS04SS /

Sample 6807 FY 2003 Project: 03-0474 Produced by: Goddard, Denise
Requestor:
Project Leader: BSTRIGGO
Facility: Gulf States Creosoting Flowood, MS
Program: SF Case No: 31635 Ending:

Media: SURFACE SOIL (0" - 12") D No: 1XY7 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

MD No: 1XY7

RESULTS	UNITS	ANALYTE		RESULTS	UNITS	ANALYTE
430 U .1	UG/KG	Benzaldehyde		430 U	UG/KG	Dibenzofuran
430 U	UG/KG	Phenol		430 U	UG/KG	2,4-Dinitrotoluene
430 U	UG/KG	bis(2-Chloroethyl) Ether		430 U	UG/KG	Diethyl Phthalate
430 U	UG/KG	2-Chlorophenol		430 U	UG/KG	Fluorene
430 U	UG/KG	2-Methylphenol	٠.,	430 U	UG/KG	4-Chlorophenyl Phenyl Ether
430 U	UG/KG .	bis(2-Chloroisopropyl) Ether		1100 U	UG/KG	4-Nitroaniline
430 U	UG/KG	Acetophenone		1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
430 U	UG/KG	(3-and/or 4-)Methylphenol		430 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
430 U	UG/KG	n-Nitrosodi-n-Propylamine		NA	UG/KG	1,2,4,5-Tetrachlorobenzene
430 U	UG/KG	Hexachloroethane		430 U	UG/KG	4-Bromophenyl Phenyl Ether
430 U	UG/KG	Nitrobenzene		430 U	UG/KG	Hexachlorobenzene (HCB)
430 U	UG/KG	Isophorone		430 U	UG/KG	Atrazine
430 U	UG/KG	2-Nitrophenol		1100 U	UG/KG	Pentachlorophenol
430 U	UG/KG	2,4-Dimethylphenol		430 U	UG/KG	Phenanthrene
430 U	·UG/KG	bis(2-Chloroethoxy)Methane		430 U	UG/KG	Anthracene
430 U	UG/KG	2,4-Dichlorophenol		430 U	UG/KG	Carbazole
430 U	UG/KG	Naphthalene		430 U	UG/KG	Di-n-Butylphthalate
430 U	UG/KG	4-Chloroaniline		430 U	UG/KG	Fluoranthene
430 U	UG/KG	Hexachlorobutadiene		430 U	UG/KG.	Pyrene
430 U	UG/KG	Caprolactam		430 UJ	UG/KG	Benzyl Butyl Phthalate
430 U	UG/KG	4-Chloro-3-Methylphenol		430 U	UG/KG	3,3'-Dichlorobenzidine
430 U	UG/KG	2-Methylnaphthalene		430 U	UG/KG	Benzo(a)Anthracene
430 U	UG/KG	Hexachlorocyclopentadiene (HCCP)		430 U	UG/KG	Chrysene
430 U	UG/KG	2,4,6-Trichlorophenol		430 UJ	UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol		430 U	UG/KG	Di-n-Octylphthalate
430 U	UG/KG	1,1-Biphenyl		430 U	UG/KG	Benzo(b)Fluoranthene
430 U	UG/KG	2-Chloronaphthalene		430 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline		430 U	UG/KG	Benzo-a-Pyrene
430 U	UG/KG	Dimethyl Phthalate		430 U	UG/KG	Indeno (1,2,3-cd) Pyrene
430 U	UG/KG	2,6-Dinitrotoluene		430 U	UG/KG	Dibenzo(a,h)Anthracene
430 U	UG/KG	Acenaphthylene		430 U	UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline		24 .	%	% Moisture
430 U	UG/KG	Acenaphthene				
1100 U	UG/KG	2,4-Dinitrophenol				
1100 U	UG/KG	4-Nitrophenol				
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U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | N-Presumptive evidence analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 06/20/2003 14:10

Sample 6807

FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/23/2003 08:30

Ending:

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS04SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

MD No: 1XY7

Inorg Contractor: SENTIN

D No: 1XY7

Org Contractor: LIBRTY

	0.11.0	-
92 J	UG/KG	Į
350 J	UG/KG	
250 J	UG/KG	2
L 0009	UG/KG	- 7

UNKNOWN ALCOHOL STIGMAST-4-EN-3-ONE 2 UNKNOWN KETONES 7 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6808 FY. 2003

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/23/2003 08:40

Ending:

**Extractables Scan** Facility: Gulf States Creosoting

ld/Station: GS04SB /

Flowood, MS

Program: SF

Case No: 31635

Inorg Contractor: SENTIN MD No: 1XY8

Media: SUBSURFACE SOIL (> 12")

Org Contractor: LIBRTY D No: 1XY8

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
430 U	UG/KG	Benzaldehyde	430 U	UG/KG	Dibenzofuran
430 U	UG/KG	Phenol	430 U	UG/KG	2,4-Dinitrotoluene
430 U	UG/KG	bis(2-Chloroethyl) Ether	430 U	UG/KG	Diethyl Phthalate
430 U	UG/KG	2-Chlorophenol	430 U	UG/KG	Fluorene
430 U	UG/KG	2-Methylphenol	430 U	UG/KG	4-Chlorophenyl Phenyl Ether
430 U	UG/KG	bis(2-Chloroisopropyl) Ether	1100 U	UG/KG	4-Nitroaniline
430 U	UG/KG	Acetophenone	1100 U	UG/KG	2-Methyl-4,6-Dinitrophenol
430 U	UG/KG	(3-and/or 4-)Methylphenol	430 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine
430 U	UG/KG	n-Nitrosodi-n-Propylamine	NA	UG/KG	1,2,4,5-Tetrachlorobenzene
430 U .	UG/KG	Hexachloroethane	430 U	UG/KG	4-Bromophenyl Phenyl Ether
430 U	UG/KG	Nitrobenzene	430 U	UG/KG	Hexachlorobenzene (HCB)
430 U	UG/KG	Isophorone	430 U	UG/KG	Atrazine
430 U	UG/KG	2-Nitrophenol	1100 U	UG/KG	Pentachlorophenol
430 U	UG/KG	2,4-Dimethylphenol	430 U	UG/KG	Phenanthrene
430 U .	UG/KG	bis(2-Chloroethoxy)Methane	430 U	UG/KG	Anthracene
430 U	UG/KG	2,4-Dichlorophenol	430 U	UG/KG	Carbazole
430 U	UG/KG	Naphthalene	430 U	UG/KG	Di-n-Butylphthalate
430 U	UG/KG	4-Chloroaniline		UG/KG	Fluoranthene
430 U	UG/KG	Hexachlorobutadiene	430 U	UG/KG	Pyrene
430 U	UG/KG	Caprolactam	430 UJ	UG/KG	Benzyl Butyl Phthalate
430 U	UG/KG	4-Chloro-3-Methylphenol	430 U	UG/KG	3,3'-Dichlorobenzidine
430 U	UG/KG	2-Methylnaphthalene	430 U	UG/KG	Benzo(a)Anthracene
430 U	UG/KG	Hexachlorocyclopentadiene (HCCP)	430 U	UG/KG	Chrysene
430 U	UG/KG	2,4,6-Trichlorophenol		UG/KG	bis(2-Ethylhexyl) Phthalate
1100 U	UG/KG	2,4,5-Trichlorophenol		UG/KG	Di-n-Octylphthalate
430 U	UG/KG	1,1-Biphenyl	430 U	UG/KG	Benzo(b)Fluoranthene
430 U	UG/KG	2-Chloronaphthalene	430 U	UG/KG	Benzo(k)Fluoranthene
1100 U	UG/KG	2-Nitroaniline	430 U	UG/KG	Benzo-a-Pyrene
430 U	UG/KG	Dimethyl Phthalate		UG/KG	Indeno (1,2,3-cd) Pyrene
430 U	UG/KG	2,6-Dinitrotoluene	430 U	UG/KG	Dibenzo(a,h)Anthracene
430 U	UG/KG	Acenaphthylene		UG/KG	Benzo(ghi)Perylene
1100 U	UG/KG	3-Nitroaniline	23	% .	% Moisture
430 U	UG/KG	Acenaphthene			
	UG/KG	2,4-Dinitrophenol			
1100 U	UG/KG	4-Nitrophenol			
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U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Project: 03-0474 Sample . 6809 FY 2003

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS05SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

MD No: 1XY9

D No: 1XY9

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/23/2003 09:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE	1			 RESULTS	UNITS	ANALYTE	•
390 U	UG/KG	Benzaldehyde				390 U	UG/KG	Dibenzofuran	
390 U	UG/KG	Phenol				390 U	UG/KG	2,4-Dinitrotoluene	
390 U	UG/KG	bis(2-Chloroethyl) Ethe	ŗ			390 U	UG/KG	Diethyl Phthalate	
390 U	UG/KG	2-Chlorophenol				390 U	UG/KG	Fluorene	
390 U	UG/KG	2-Methylphenol				390 U	UG/KG	4-Chlorophenyl Phenyl Ether	•
390 U	UG/KG	bis(2-Chloroisopropyl)	ther			980 U	UG/KG	4-Nitroaniline	
390 U	UG/KG	Acetophenone				980 U	UG/KG	2-Methyl-4,6-Dinitrophenol	
390 U	UG/KG	(3-and/or 4-)Methylpher	iol			390 U	UG/KG	n-Nitrosodiphenylamine/Diphenylamine	•
390 U	UG/KG	n-Nitrosodi-n-Propylami	ine			NĄ	UG/KG	1,2,4,5-Tetrachlorobenzene	
390 U	UG/KG	Hexachloroethane				390 U	UG/KG	4-Bromophenyl Phenyl Ether	•
390 U	UG/KG	Nitrobenzene		•		、390 U	UG/KG .	Hexachlorobenzene (HCB)	
390 U	UG/KG	Isophorone		•	_	390 U	UG/KG	Atrazine	•
390 U	UG/KG	2-Nitrophenol	1	•		980 U	UG/KG	Pentachlorophenol	•
390 U	UG/KG	2,4-Dimethylphenol		•	•	320 J	UG/KG	Phenanthrene	:
390 U	UG/KG	bis(2-Chloroethoxy)Met	hane			470	UG/KG	Anthracene	•
390 U	UG/KG	2,4-Dichlorophenol	1			200 J	UG/KG	Carbazole	
390 U	UG/KG	Naphthalene	*			390 U	UG/ĶG	Di-n-Butylphthalate	
390 U	UG/KG	4-Chloroaniline	1			2600	UG/KG	Fluoranthene	
390 U	UG/KG	Hexachlorobutadiene	İ		•	1500	UG/KG	Pyrene	•
390 U	UG/KG	Caprolactam		-		390 UJ	UG/KG	Benzyl Butyl Phthalate	
390 U	UG/KG	4-Chloro-3-Methylpheno	ήl ·			390 U	UG/KG	3,3'-Dichlorobenzidine	·
390 U	UG/KG	2-Methylnaphthalene		•		770	UG/KG	Benzo(a)Anthracene	
390 U	UG/KG	Hexachlorocyclopentad	iene (HCCP)			1500	UG/KG	Chrysene	
390 U	UG/KG	2,4,6-Trichlorophenol				390 UJ	UG/KG	bis(2-Ethylhexyl) Phthalate	
980 U	UG/KG	2,4,5-Trichlorophenol			, ,	390 U	UG/KG ·	Di-n-Octylphthalate	
390 ⊍	UG/KG	1,1-Biphenyl	į	•		1800 J	UG/KG	Benzo(b)Fluoranthene	
390 U	UG/KG	2-Chloronaphthalene	T.			2100 J	UG/KG	Benzo(k)Fluoranthene	
	UG/KG	2-Nitroaniline	1	•		810	UG/KG	Benzo-a-Pyrene	•
	UG/KG	Dimethyl Phthalate				610	UG/KG	Indeno (1,2,3-cd) Pyrene	
390 U	UG/KG	2,6-Dinitrotoluene		-		190 J	UG/KG	Dibenzo(a,h)Anthracene	
	UG/KG	Acenaphthylene	İ	•		380 J	UG/KG	Benzo(ghi)Perylene	
	UG/KG	3-Nitroaniline	1	•		15	%	% Moisture	
	UG/KG	Acenaphthene						·	
	UG/KG	2,4-Dinitrophenol						•	
980 U	UG/KG	4-Nitrophenol							
			:					•	
			!				-	•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Production Date: 06/20/2003 14:10

Sample

6809

FY 2003

Project: 03-0474

**MISCELLANEOUS COMPOUNDS** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS05SS /

MD No: 1XY9 D No: 1XY9

Media: SURFACE SOIL (0" - 12")

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:15

Ending:

RESULTS	UNITS	ANALYTE
190 J	UG/KG	METHYLANTHRACENE
260 NJ	UG/KG	9, 10-ANTHRACENEDIONE
160 NJ	UG/KG	PHENANTHRENE, 2,3-DIMETHYL-
160 J	UG/KG	BENZOFLUORENE
470 NJ	UG/KG	11H-BENZO [A] FLUORENE
250 NJ	UG/KG	PYRENE, 2-METHYL-
280 NJ	UG/KG	PYRENE, 1,3-DIMETHYL-
380 NJ	UG/KG	BENZO [B] NAPHTHO [2,1-D] THIOPHENE
310 NJ	UG/KG	7H-BENZ [DE] ANTHRACEN-7-ONE
270 NJ	UG/KG	CHRYSENE, 6-METHYL-
330 NJ	UG/KG	5,12-NAPHTHACENEDIONE
290 NJ	UG/KG	BENZO [E] PYRENE
260 NJ	UG/KG	1,2:7,8-DIBENZPHENANTHRENE
180 NJ	UG/KG	[3,4:9,10] DIBENZPYRENE
1500 J	UG/KG	4 UNKNOWN PAHS
2700 J	UG/KG	11 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

6810 FY 2003 Project: 03-0474 Sample

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:25

Ending:

Flowood, MS Facility: Gulf States Creosoting

Program: SF

Case No: 31635

Id/Station: GS05SB /

**Extractables Scan** 

MD No: 1XZ0

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

Org Contractor: LIBRTY D No: 1XZ0

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE
410 U	UG/KG	Benzaldehyde	1		410 U	UG/KG	Dibenzofuran
410 U	UG/KG	Phenol			410 U	UG/KG	2.4-Dinitrotoluene
410 U	UG/KG	bis(2-Chloroethyl) Ether		•	410 U	UG/KG	Diethyl Phthalate
410 U	UG/KG	2-Chiorophenoi			410 U	UG/KG	Fluorene
410 U	UG/KG	2-Methylphenol	·		410 U	UG/KG	4-Chlorophenyl Phenyl Ether
410 U	UG/KG	bis(2-Chloroisopropyl) E	: Ethor		1000 U	UG/KG	4-Onlorophenyl Frienyl Ethel
410 U	UG/KG	Acetophenone	iner		1000 U	UG/KG	2-Methyl-4,6-Dinitrophenol
	UG/KG UG/KG	(3-and/or 4-)Methylpher			410 U	UG/KG	
410 U					410 U NA		n-Nitrosodiphenylamine/Diphenylamine
410 U	UG/KG	n-Nitrosodi-n-Propylami Hexachloroethane				UG/KG	1,2,4,5-Tetrachlorobenzene
410 U	UG/KG				410 U		4-Bromophenyl Phenyl Ether
410 U ·	UG/KG	Nitrobenzene			410 U	UG/KG	Hexachlorobenzene (HCB)
410 U	UG/KG	Isophorone			410 U	UG/KG	Atrazine
410 U	UG/KG	2-Nitrophenol			1000 U	UG/KG	Pentachlorophenol
410 U	UG/KG	2,4-Dimethylphenol			410 U	UG/KG	Phenanthrene
410 U	UG/KG	bis(2-Chloroethoxy)Met	nane		410 U	UG/KG	Anthracene
410 U	UG/KG	2,4-Dichlorophenol			410 U	UG/KG	Carbazole
410 U	UG/KG	Naphthalene	_		410 U	UG/KG	Di-n-Butylphthalate
410 U	UG/KG	4-Chloroaniline			410 U	UG/KG	Fluoranthene
410 U	UG/KG	Hexachlorobutadiene			410 U	UG/KG	Pyrene
410 U	UG/KG	Caprolactam			410 UJ	UG/KG	Benzyl Butyl Phthalate
410 U	UG/KG	4-Chloro-3-Methylpheno	א 		410 Ú	UG/KG	3,3'-Dichlorobenzidine
410 U	UG/KG UG/KG	2-Methylnaphthalene	(UCCD)		410 U	UG/KG	Benzo(a)Anthracene
410 U		Hexachlorocyclopentadi	ene (nccr)		410 U	UG/KG	Chrysene
410 U	UG/KG	2,4,6-Trichlorophenol			450 J	UG/KG	bis(2-Ethylhexyl) Phthalate
1000 U	UG/KG	2,4,5-Trichlorophenol			410 U	UG/KG	Di-n-Octylphthalate
410 U	UG/KG	1,1-Biphenyl			410 U	UG/KG	Benzo(b)Fluoranthene
410 U	UG/KG	2-Chloronaphthalene			410 U	UG/KG	Benzo(k)Fluoranthene
1000 U	UG/KG	2-Nitroaniline			410 U	UG/KG	Benzo-a-Pyrene
410 U	UG/KG	Dimethyl Phthalate	l· I		410 U	UG/KG	Indeno (1,2,3-cd) Pyrene
410 U	UG/KG	2,6-Dinitrotoluene			410 U	UG/KG	Dibenzo(a,h)Anthracene
410 U	UG/KG	Acenaphthylene			410 U	UG/KG	Benzo(ghi)Perylene
1000 U	UG/KG	3-Nitroaniline			19	%	% Moisture
410 U	UG/KG	Acenaphthene					
1000 U	UG/KG	2,4-Dinitrophenol	,			7	
1000 U	UG/KG	4-Nitrophenol					
•			ļ				

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6811 FY 2003 Project: 03-0474

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS03GW /

Case No: 31635

Media: GROUNDWATER

MD No: 1XZ1

D No: 1XZ1

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:55

Ending:

RESULTS	UNITS	ANALYTE					RESULTS	UNITS	ANALYTE	
5.0 U	UG/L	Benzaldehyde				•	5.0 U	UG/L	Dibenzofuran	
5.0 U •	UG/L	Phenol				-	5.0 U	UG/L	2.4-Dinitrotoluene	•
5.0 U	UG/L	bis(2-Chloroethyl) Ethe	; <b>r</b>	• • •		•	5.0 UJ	UG/L	Diethyl Phthalate	•
5.0 U	UG/L	2-Chlorophenol					5.0 U	UG/L	Fluorene	
5.0 U	UG/L	2-Methylphenol	· ·	•	•		5.0 U	UG/L	4-Chlorophenyl Phenyl Ether	
5.0 U	UG/L	bis(2-Chloroisopropyl) I	Ether		•		20 U	UG/L	4-Nitroaniline	
5.0 U	UG/L	Acetophenone	!				20 UJ	UG/L	2-Methyl-4,6-Dinitrophenol	
. 5.0 U	UG/L	(3-and/or 4-)Methylphei	hol		•		5.0 U	UG/L	n-Nitrosodiphenylamine/Diphenylan	nine
5.0 U	UG/L	n-Nitrosodi-n-Propylam				· · ·	5.0 U	UG/L	1,2,4,5-Tetrachlorobenzene	
5.0 U	UG/L	Hexachloroethane					5.0 U	UG/L	4-Bromophenyl Phenyl Ether	
5.0 U	UG/L	Nitrobenzene					5.0 U	UG/L	Hexachlorobenzene (HCB)	
5.0 U	UG/L	Isophorone					5.0 U	UG/L	Atrazine	
5.0 U	UG/L	2-Nitrophenol					5.0 U	UG/L	Pentachlorophenol	
5.0 U	UG/L	2,4-Dimethylphenol	ļ			• •	5.0 U	UG/L	Phenanthrene	•
5.0 U	UG/L	bis(2-Chloroethoxy)Met	hane				5.0 U	UG/L	Anthracene	
5.0 U	UG/L	2,4-Dichlorophenol	£				· NA	UG/L	Carbazole	
5.0 U	UG/L	Naphthalene	•				5.0 UJ	UG/L	Di-n-Butylphthalate	
5.0 U	UG/L	4-Chloroaniline		-	-		5.0 U	UG/L	Fluoranthene	
5.0 U	UG/L	Hexachlorobutadiene					5.0 U	UG/L	Pyrene	
5.0 UJ	UG/L	Caprolactam	į				5.0 UJ	UG/L	Benzyl Butyl Phthalate	
5.0 U	UG/L	4-Chloro-3-Methylpheno	ρl	•			5.0 UJ	UG/L	3,3'-Dichlorobenzidine	
5.0 U	UG/L	2-Methylnaphthalene	!				5.0 U	UG/L	Benzo(a)Anthracene	
5.0 UJ	UG/L	Hexachlorocyclopentad	iene (HCCP)				5.0 U	UG/L	Chrysene	
5.0 U	UG/L	2,4,6-Trichlorophenol	1		•		5.0 UJ	UG/L	bis(2-Ethylhexyl) Phthalate	
20 U	UG/L	2,4,5-Trichlorophenol			• •		5.0 UJ	UG/L	Di-n-Octylphthalate	
5.0 UJ	UG/L	1,1-Biphenyl					5.0 U	UG/L	Benzo(b)Fluoranthene	
5.0 U	UG/L	2-Chioronaphthalene	1				5.0 U	UG/L	Benzo(k)Fluoranthene	
20 U	UG/L	2-Nitroaniline	<u>.</u>				5.0 U	UG/L	Benzo-a-Pyrene	•
	UG/L	Dimethyl Phthalate					5.0 U	UG/L	Indeno (1,2,3-cd) Pyrene	•
	UG/L	2,6-Dinitrotoluene	: i	•			- 5.0 U	UG/L	Dibenzo(a,h)Anthracene	
	UG/L	Acenaphthylene				•	5.0 U	UG/L	Benzo(ghi)Perylene	
	UG/L	3-Nitroaniline				•				
	UG/L	Acenaphthene	!							
	UG/L	2,4-Dinitrophenol							•	
20 U	UG/L	4-Nitrophenol								
			1						·	
	•				•					•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Produced by: Goddard, Denise

Sample 6811 FY 2003 Project: 03-0474

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS03GW /

MD No: 1XZ1

D No: 1XZ1

Inorg Contractor: SENTIN

Org Contractor: A4

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:55

Ending:

**RESULTS UNITS ANALYTE** 2.1 NJ UG/L LIMONENE

Media: GROUNDWATER

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/23/2003 13:20

Requestor:

Ending:

Sample 6812 FY 2003 Project: 03-0474

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS04GW /

Media: GROUNDWATER

MD No: 1XZ3

Inorg Contractor: SENTIN

D No: 1XZ3

Org Contractor: A4

RESULTS	UNITS	ANALYTE			RESULTS	UNITS	ANALYTE
5.0 U	UG/L	Benzaldehyde			5.0 U	UG/L	Dibenzofuran
5.0 U	ÜG/L	Phenol			5.0 U	UG/L	2,4-Dinitrotoluene
5.0 U	UG/L	bis(2-Chloroethyl) Ether	•		5.0 U	UG/L	Diethyl Phthalate
5.0 U	UG/L	2-Chlorophenol		•	5.0 U	UG/L	Fluorene
5.0 U	UG/L	2-Methylphenol			5.0 U	UG/L	4-Chlorophenyl Phenyl Ether
- ∕5.0 U	UG/L	bis(2-Chloroisopropyl) Ether		-	20 U	UG/L	4-Nitroaniline
5.0 U	UG/L	Acetophenone		•	20 UJ	UG/L	2-Methyl-4,6-Dinitrophenol
5.0 U	UG/L	(3-and/or 4-)Methylphenol	•		5.0 U	UG/L	n-Nitrosodiphenylamine/Diphenylamine
5.0 U	UG/L	n-Nitrosodi-n-Propylamine		. •	5.0 U	UG/L	1,2,4,5-Tetrachlorobenzene
5.0 U	UG/L	Hexachloroethane			5.0 U	UG/L	4-Bromophenyl Phenyl Ether
5.0 U	UG/L	Nitrobenzene			5.0 U	UG/L	Hexachlorobenzene (HCB)
5.0 U	UG/L	Isophorone			5.0 U	UG/L	Atrazine
5.0 U	UG/L	2-Nitrophenol			5.0 U	UG/L	Pentachlorophenol
5.0 U	UG/L	2,4-Dimethylphenol			5.0 U	UG/L	Phenanthrene
5.0 U	UG/L	bis(2-Chloroethoxy)Methane		•	5.0 U	UG/L	Anthracene
5.0 U	UG/L	2,4-Dichlorophenol	•		NA	UG/L	Carbazole
5.0 U	UG/L	Naphthalene	•		5.0 U	UG/L	Di-n-Butylphthalate
5.0 U	UG/L	4-Chloroaniline			5.0 U	UG/L	Fluoranthene
5.0 U	UG/L	Hexachlorobutadiene			5.0 U	UG/L	Pyrene
5.0 U	UG/L	Caprolactam	•		5.0 U	UG/L	Benzyl Butyl Phthalate
5.0 U	UG/L	4-Chloro-3-Methylphenol			5.0 UJ	UG/L	3,3'-Dichlorobenzidine
5.0 U	UG/L	2-Methylnaphthalene		-	5.0 U	UG/L	Benzo(a)Anthracene
5.0 UJ	UG/L:	Hexachlorocyclopentadiene (HCCP)			5.0 U	UG/L	Chrysene
5.0 U	UG/L	2,4,6-Trichlorophenol			5.0 U	UG/L	bis(2-Ethylhexyl) Phthalate
20 Ü	UG/L	2,4,5-Trichlorophenol			5.0 U	UG/L	Di-n-Octylphthalate
5.0 U	UG/L	1,1-Biphenyl			5.0 U	UG/L	Benzo(b)Fluoranthene
5.0 U	UG/L	2-Chloronaphthalene	•		5.0 U	UG/L	Benzo(k)Fluoranthene
	UG/L	2-Nitroaniline	•		5.0 U	UG/L	Benzo-a-Pyrene
5.0 U	UG/L	Dimethyl Phthalate	•		5.0 U	UG/L	Indeno (1,2,3-cd) Pyrene
	UG/L	2,6-Dinitrotoluene			5.0 U	UG/L	Dibenzo(a,h)Anthracene
5.0 U	UG/L	Acenaphthylene			5.0 U	UG/L	Benzo(ghi)Perylene
20 U	UG/L	3-Nitroaniline		•			
	UG/L	Acenaphthene					
	UG/L	2,4-Dinitrophenol	•				•
20 U	UG/L	4-Nitrophenol					
		;					

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Production Date: 06/20/2003 14:10

Sample

6813 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/23/2003 11:45

Ending:

**Extractables Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS05GW /

MD No: 1XZ4

Inorg Contractor: SENTIN

Media: GROUNDWATER D No: 1XZ4

Org Contractor: A4

R	ESULTS	UNITS	ANALYTE	RESULTS	UNITS	ANALYTE
	5.0 U	UG/L	Benzaldehyde	5.0 U	UG/L	Dibenzofuran
	5.0 U	UG/L	Phenol	5.0 U	UG/L	2,4-Dinitrotoluene
	5.0 U *	UG/L	bis(2-Chloroethyl) Ether	5.0 U	UG/L	Diethyl Phthalate
	5.0 U	UG/L	2-Chlorophenol	5.0 U	UG/L	Fluorene
	5.0 U	UG/L	2-Methylphenol	5.0 U	UG/L	4-Chlorophenyl Phenyl Ether
	5.0 U	UG/L	bis(2-Chloroisopropyl) Ether	20 U	UG/L	4-Nitroaniline
	5.0 U	UG/L	Acetophenone	20 UJ	UG/L	2-Methyl-4,6-Dinitrophenol
	5.0 U	UG/L	(3-and/or 4-)Methylphenol	5.0 U	UG/L	n-Nitrosodiphenylamine/Diphenylamine
		UG/L	n-Nitrosodi-n-Propylamine	5.0 U	UG/L	1,2,4,5-Tetrachlorobenzene
	5.0 U	UG/L	Hexachloroethane	5.0 U	UG/L	4-Bromophenyl Phenyl Ether
		UG/L	Nitrobenzene	5.0 U	UG/L	Hexachlorobenzene (HCB)
	·5.0 U	UG/L	Isophorone	5.0 U	UG/L	Atrazine
		UG/L	2-Nitrophenol	5.0 U	UG/L	Pentachlorophenol
		UG/L	2,4-Dimethylphenol	5.0 U	UG/L	Phenanthrene
	5.0 U	ÜG/L	bis(2-Chloroethoxy)Methane	5.0 U	UG/L	Anthracene
	5.0 U	UG/L	2,4-Dichlorophenol	NA	UG/L	Carbazole
	5.0 U	UG/L	Naphthalene	5.0 U	UG/L	Di-n-Butylphthalate
		UG/L	4-Chloroaniline	5.0 U	UG/L	Fluoranthene
		UG/L	Hexachlorobutadiene	5.0 U	UG/L	Pyrene
		UG/L	Caprolactam	5.0 U	UG/L	Benzyl Butyl Phthalate
		UG/L	4-Chloro-3-Methylphenol	5.0 UJ	UG/L	3,3'-Dichlorobenzidine
		UG/Ļ	2-Methylnaphthalene	5.0 U	UG/L	Benzo(a)Anthracene
		UG/Ĺ	Hexachlorocyclopentadiene (HCCP)	5.0 U	UG/L	Chrysene
		UG/L	2,4,6-Trichlorophenol	5.0 U	UG/L	bis(2-Ethylhexyl) Phthalate
		UG/L	2,4,5-Trichlorophenol	5.0 U	UG/L	Di-n-Octylphthalate
		UG/L	1,1-Biphenyl	5.0 U	UG/L	Benzo(b)Fluoranthene
		UG/L	2-Chloronaphthalene	5.0 U	UG/L	Benzo(k)Fluoranthene
		UG/L	2-Nitroaniline	5.0 U	UG/L	Benzo-a-Pyrene
		UG/L	Dimethyl Phthalate	5.0 U	UG/L	Indeno (1,2,3-cd) Pyrene
		UG/L	2,6-Dinitrotoluene	5.0 U	UG/L	Dibenzo(a,h)Anthracene
		UG/L	Acenaphthylene	5.0 U	UG/L	Benzo(ghi)Perylene
		UG/L	3-Nitroaniline			•
			Acenaphthene			
			2,4-Dinitrophenol			
	20 U	UG/L	4-Nitrophenol			
						•
			· i			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EXTRACTABLES SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:10

Sample

6813 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 11:45

Ending:

MISCELLANEOUS COMPOUNDS

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS05GW /

Case No: 31635

MD No: 1XZ4

Inorg Contractor: SENTIN

Media: GROUNDWATER

D No: 1XZ4

Org Contractor: A4

**RESULTS UNITS** 

ANALYTE-

4.7 J UG/L 2 UNKNOWN COMPOUNDS

Data Reported as Identified by CLP Lab - IDs Not Verified

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.



# UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

# Region 4

Science and Ecosystem Support Division 980 College Station Road Athens, Georgia 30605-2720

## **MEMORANDUM**

Date: 06/20/2003

Subject: Results of PESTICIDES/PCB Sample Analysis

03-0474

**Gulf States Creosoting** 

Flowood, MS

From: Goddard, Denise

To: Striggow, Brian

Thru: QA Office.

Attached are the results of analysis of samples collected as part of the subject project. If you have any questions, please contact me.

FY 2003 Project: 03-0474 Sample 6754

Produced by: Goddard, Denise

Beginning: 04/22/2003 09:00

**Pesticides Scan** 

Requestor: Project Leader: BSTRIGGO

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS01SD / MD No: 1XR4 Media: SEDIMENT D No: 1XR4

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
3.8 U	UG/KG	alpha-BHC
3.8 U	UG/KG	beta-BHC
3.8 U .	UG/KG	delta-BHC
3.8 U	UG/KG	gamma-BHC (Lindane)
3.8 U	UG/KG	Heptachlor
3.8 U	UG/KG	Aldrin
3.8 U	UG/KG	Heptachlor Epoxide
3.8 U	UG/KG	Endosulfan I (alpha)
7.3 U	UG/KG	Dieldrin
7.3 U	UG/KG	4,4'-DDE (p,p'-DDE)
7.3 U	UG/KG	Endrin
7.3 U	UG/KG	Endosulfan II (beta)
7.3 U	UG/KG -	4,4'-DDD (p,p'-DDD)
7.3 U	UG/KG	Endosulfan Sulfate
7.3 U	UG/KG	4,4'-DDT (p,p'-DDT)
38 U	UG/KG	Methoxychlor
7.3·U	UG/KG	Endrin Ketone
7.3 U	UG/KG	Endrin Aldehyde
3.8 U	UG/KG	alpha-Chlordane /2
3.8 U	UG/KG	gamma-Chlordane /2
380 U <sub>.</sub>	UG/KG	Toxaphene
73 U	UG/KG	PCB-1016 (Aroclor 1016)
150 U	UG/KG	PCB-1221 (Aroclor 1221)
73 U	UG/KG	PCB-1232 (Aroclor 1232)
73 U	UG/KG	PCB-1242 (Aroclor 1242)
73 U	UG/KG	PCB-1248 (Aroclor 1248)
73 U	UG/KG	PCB-1254 (Aroclor 1254)
73 U	UG/KG	PCB-1260 (Aroclor 1260)
55	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

PESTICIDES	/PCB	SAMPLE	ANALYSIS

Production Date: 06/20/2003 14:11

Project: 03-0474 Sample 6755 FY 2003

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF Id/Station: GS18SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XR5

D No: 1XR5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:50

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.3 U	UG/KG	alpha-BHC
2.3 U	UG/KG	beta-BHC
2.3 U	UG/KG	delta-BHC
1.7 J	UG/KG	gamma-BHC (Lindane)
2.3 U	UG/KG	Heptachlor
2.3 U	UG/KG	Aldrin
2.3 U	UG/KG	Heptachlor Epoxide
2.3 U	UG/KG	Endosulfan I (alpha)
4.4 U	UG/KG	Dieldrin
4.4 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.4 U	UG/KG	Endrin
4.4 U	UG/KG	Endosulfan II (beta)
4.4 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.4 U	UG/KG	Endosulfan Sulfate
4.4 U	UG/KG	4,4'-DDT (p,p'-DDT)
23 U	UG/KG	Methoxychlor
4.4 U	UG/KG	Endrin Ketone
4.4 U	UG/KG	Endrin Aldehyde
2.3 U	UG/KG	alpha-Chlordane /2
2.3 U	UG/KG	gamma-Chlordane /2
230 U	UG/KG	Toxaphene
44 U	UG/KG	PCB-1016 (Aroclor 1016)
89 U	UG/KG	PCB-1221 (Aroclor 1221)
44 U	UG/KG	PCB-1232 (Aroclor 1232)
44 U	UG/KG	PCB-1242 (Aroclor 1242)
44 U	UG/KG	PCB-1248 (Aroclor 1248)
44 U	UG/KG	PCB-1254 (Aroclor 1254)
42 J	UG/KG	PCB-1260 (Aroclor 1260)
25	%	% Moisture
		į

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

PESTICIDES/PCB	CAMPIE	: ANAI	VCIC
PESTICIDES/FCD	SAINFLE	. AIXAL	_ 1 - 31 - 3

Production Date: 06/20/2003 14:11

Sample 6756 FY 2003 Project: 03-0474

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS18SS /

Media: SURFACE SOIL (0" - 12")

Case No: 31635

MD No: 1XR6

D No: 1XR6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.3 U	UG/KG	Dieldrin
4.3 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.3 U	UG/KG	Endrin
4.3 U	UG/KG	Endosulfan II (beta)
4.3 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.3 U	UG/KG	Endosulfan Sulfate
4.3 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.3 U	UG/KG	Endrin Ketone
4.3 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U	UG/KG	Toxaphene
43 U	UG/KG	PCB-1016 (Aroclor 1016)
87 U	UG/KG	PCB-1221 (Aroclor 1221)
43 U	UG/KG	PCB-1232 (Aroclor 1232)
43 U	UG/KG	PCB-1242 (Aroclor 1242)
43 U	UG/KG	PCB-1248 (Aroclor 1248)
43 U	UG/KG	PCB-1254 (Aroclor 1254)
43 U	UG/KG	PCB-1260 (Aroclor 1260)
23	%	% Moisture
		j

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 6757 FY 2003 Sample

Pesticides Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS19SB /

Media: SUBSURFACE SOIL (> 12")

MD No: 1XR7

D No: 1XR7

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:35

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.2 U	UG/KG	Dieldrin
4.2 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.2 U	UG/KG	Endrin
4.2 U	UG/KG	Endosulfan II (beta)
4.2 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.2 U	UG/KG	Endosulfan Sulfate
4.2 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.2 U	UG/KG	Endrin Ketone
4.2 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U	UG/KG	Toxaphene
42 U	UG/KG	PCB-1016 (Arocior 1016)
85 U -	UG/KG	PCB-1221 (Aroclor 1221)
42 U	UG/KG	PCB-1232 (Aroclor 1232)
42 U	UG/KG	PCB-1242 (Aroclor 1242)
42 U	UG/KG	PCB-1248 (Aroclor 1248)
42 U	UG/KG	PCB-1254 (Aroclor 1254)
. 39 J	UG/KG	PCB-1260 (Aroclor 1260)
<sub>.</sub> 21	%	% Moisture
		; (

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample FY 2003 Project: 03-0474 6758

Flowood, MS

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 09:30

Ending:

Case No: 31635

Media: SURFACE SOIL (0" - 12")

Facility: Gulf States Creosoting

**Pesticides Scan** 

Id/Station: GS19SS /

Program: SF

MD No: 1XR8 D No: 1XR8

Inorg Contractor: SENTIN Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
2.0 Ú	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.0 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
4.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
20 U	UG/KG	Methoxychior
4.0 U	UG/KG	Endrin Ketone
4.0 U	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
	UG/KG	PCB-1016 (Aroclor 1016)
81 U	ÚG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
40 U	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
, 17	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value,

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474 Sample FY 2003

Pesticides Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS02SD /

Case No: 31635

Media: SEDIMENT

MD No: 1XR9

D No: 1XR9

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:50

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
3.5 U	UG/KG	alpha-BHC
3.5 U	UG/KG	beta-BHC
3.5 U	UG/KG	delta-BHC
3.5 U	UG/KG	gamma-BHC (Lindane)
3.5 U	UG/KG	Heptachlor
3.5 U	UG/KG	Aldrin
3.5 U	UG/KG	Heptachlor Epoxide
3.5 U	UG/KG	Endosulfan I (alpha)
6.9 U	UG/KG	Dieldrin
7.9	UG/KG	4,4'-DDE (p,p'-DDE)
6.9 U	UG/KG	Endrin
6.5 J	UG/KG	Endosulfan II (beta)
6.9 U	UG/KG	4,4'-DDD (p,p'-DDD)
10 N	UG/KG	Endosulfan Sulfate
6.9 U	UG/KG	4,4'-DDT (p,p'-DDT)
35 U	UG/KG	Methoxychlor
6.9 U	UG/KG	Endrin Ketone
6.9 U	UG/KG	Endrin Aldehyde
3.5 U	UG/KG	alpha-Chlordane /2
3.5 U	UG/KG	gamma-Chlordane /2
350 U	UG/KG	Toxaphene
69 U	UG/KG	PCB-1016 (Aroclor 1016)
140 U	UG/KG	PCB-1221 (Aroclor 1221)
69 U	UG/KG	PCB-1232 (Aroclor 1232)
69 U	UG/KG	PCB-1242 (Aroclor 1242)
69 U	UG/KG	PCB-1248 (Aroclor 1248)
69 U	UG/KG	PCB-1254 (Aroclor 1254)
69 U	UG/KG	PCB-1260 (Aroclor 1260)
52	%	% Moisture
		:

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

DECTIONEC/DOD	CARADIE	ANIAL VOIC
PESTICIDES/PCB	SAMPLE	ANALYSIS

Production Date: 06/20/2003 14:11

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise Sample 6760 FY 2003 Project: 03-0474 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 10:05 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS17SS / MD No: 1XS0 Org Contractor: LIBRTY D No: 1XS0 Media: SURFACE SOIL (0" - 12")

	<u> </u>		
RESULTS UNITS	ANALYTE		
2.2 U UG/KG	alpha-BHC		·
2.2 U UG/KG	beta-BHC		•
2.2 U UG/KG	delta-BHC	•	
2.2 U . UG/KG	gamma-BHC (Lindane)		
2.2 U UG/KG	Heptachlor		
2.2 U UG/KG	Aldrin		
2.2 U UG/KG	Heptachlor Epoxide		_
2.2 U UG/KG	Endosulfan I (alpha)		•
4.2 U UG/KG	Dieldrin	•	
4.2 U UG/KG	4,4'-DDE (p,p'-DDE)		•
6.4 J UG/KG	Endrin	•	
4.2 U UG/KG	Endosulfan II (beta)		
4.2 U UG/KG	4,4'-DDD (p,p'-DDD)		•
4.2 U UG/KG	Endosulfan Sulfate		
4.2 U UG/KG	4,4'-DDT (p,p'-DDT)		•
22 U UG/KG	Methoxychlor		
20 J UG/KG	Endrin Ketone		
4.2 U UG/KG	Endrin Aldehyde		
2.2 U UG/KG	alpha-Chlordane /2	•	
2.2 U UG/KG	gamma-Chlordane /2		
220 Ü UG/KG	Toxaphene		
42 U UG/KG	PCB-1016 (Aroclor 1016)		
85 U UG/KG	PCB-1221 (Aroclor 1221)		
42 U UG/KG	PCB-1232 (Aroclor 1232)		
42 U UG/KG	PCB-1242 (Aroclor 1242)		
42 U UG/KG	PCB-1248 (Aroclor 1248)	•	
42 U UG/KG	PCB-1254 (Aroclor 1254)	* .	
42 U UG/KG	PCB-1260 (Aroclor 1260)		
21 %	% Moisture		
	:		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

PESTICIDES/PCB	SAMPLE	<b>ANALYSIS</b>

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6761 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Flawood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 10:20 Program: SF Case No: 31635 Ending: inorg Contractor: SENTIN Id/Station: GS17SB / MD No: 1XS1 D No: 1XS1 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") DATA REPORTED ON DRY WEIGHT BASIS

		, , , , , , , , , , , , , , , , , , , ,		<b>3</b>	DATA NEI ONTED ON DITT WEIGHT BA
RESULTS	UNITS	ANALYTE			
2.2 U	UG/KG	alpha-BHC			
2.2 U	UG/KG	beta-BHC		•	
2.2 U	UG/KG	delta-BHC			·
2.2 U	UG/KG	gamma-BHC (Lindane)	·		
2.2 U	UG/KG	Heptachlor		•	
2.2 U	UG/KG	Aldrin	•		
2.2 U	UG/KG	Heptachlor Epoxide	•		
2.2 U	UG/KG	Endosulfan I (alpha)			·
4.2 U	UG/KG	Dieldrin			
4.2 U	UG/KG	4,4'-DDE (p,p'-DDE)			
4.2 U	UG/KG	Endrin	•		•
4.2 U	UG/KG	Endosulfan II (beta)		•	•
.4.2 U	UG/KG	4,4'-DDD (p,p'-DDD)	•	•	•
4.2 U	UG/KG	Endosulfan Sulfate	•		
4.2 U	UG/KG	4,4'-DDT (p,p'-DDT)			•
22 U	UG/KG	Methoxychlor			
4.2 U	UG/KG	Endrin Ketone		•	•
4.2 U	UG/KG	Endrin Aldehyde			•
.2.2 U	UG/KG	alpha-Chlordane /2		•	•
2.2 U	UG/KG	gamma-Chlordane /2			
220 U	UG/KG	Toxaphene		•	
42 U	UG/KG	PCB-1016 (Aroclor 1016			
86 U -	UG/KG	PCB-1221 (Aroclor 1221			
42 U	UG/KG	PCB-1232 (Aroclor 1232			
42 U	UG/KG	PCB-1242 (Aroclor 1242)		•	
42 U	UG/KG	PCB-1248 (Aroclor 1248			
42 U	UG/KG	PCB-1254 (Aroclor 1254			
42 U	UG/KG	PCB-1260 (Aroclor 1260	)		
22	%	% Moisture			
				•	•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

#### PESTICIDES/PCB SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:11

Sample 6762 FY 2003 Project: 03-0474

Facility: Gulf States Creosoting

Program: SF

**Pesticides Scan** 

Id/Station: GS20SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635

MD No: 1XS2

D No: 1XS2

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:45

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KĠ	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.2 U	UG/KG	Dieldrin
4.1 NJ	UG/KG	4,4'-DDE (p,p'-DDE)
4.2 U	UG/KG	Endrin
4.2 U	UG/KG	Endosulfan II (beta)
4.2 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.2 U -	UG/KG	Endosulfan Sulfate
4.2 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.2 U	UG/KG	Endrin Ketone
4.2 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U	UG/KG	Toxaphene
42 U	UG/KG	PCB-1016 (Aroclor 1016)
85 U	UG/KG	PCB-1221 (Aroclor 1221)
42 U	UG/KG	PCB-1232 (Aroclor 1232)
42 U	UG/KG	PCB-1242 (Aroclor 1242)
42 U .	UG/KG	PCB-1248 (Aroclor 1248)
42 U	UG/KG	PCB-1254 (Aroclor 1254)
42 U	UG/KG	PCB-1260 (Aroclor 1260)
21	%	% Moisture
		<u> </u>

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6763 FY 2003

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS20SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

MD No: 1XS3

D No: 1XS3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:55

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.1 U	UG/KG	alpha-BHC
2.1 U	UG/KG	beta-BHC
2.1 U	UG/KG	delta-BHC
2.1 U	UG/KG	gamma-BHC (Lindane)
2.1 U •	UG/KG	Heptachlor
2.1 U	UG/KG	Aldrin
2.1 U	UG/KG	Heptachlor Epoxide
2.1 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.0 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
4.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
21 U	UG/KG	Methoxychlor
4.0 U	UG/KG	Endrin Ketone
4.0 U	UG/KG	Endrin Aldehyde
2.1 U	UG/KG	alpha-Chlordane /2
2.1 U	UG/KG	gamma-Chlordane /2
210 U	UG/KG	Toxaphene
40 U	UG/KG	PCB-1016 (Aroclor 1016)
82 U	UG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
18	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

PESTI	CIDES	/PCB	SAMPL	.E A	NAL	YSIS

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6764 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 10:45 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS20SD / MD No: 1XS4 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XS4 DATA REPORTED ON DRY WEIGHT BASIS

		· · · · · · · · · · · · · · · · · · ·
RESULTS	UNITS	ANALYTE
2.1 U	UG/KG	alpha-BHC
2.1 U	UG/KG	beta-BHC
2.1 U	UG/KG	delta-BHC
2.1 U	UG/KG	gamma-BHC (Lindane)
2.1 U	UG/KG	Heptachlor
2.1 U	UG/KG	Aldrin
2.1 U	UG/KG	Heptachlor Epoxide
2.1 U	UG/KG	Endosulfan I (alpha)
4.1 U	UG/KG	Dieldrin
3.3 NJ	UG/KG	4,4'-DDE (p,p'-DDE)
4.1 U	UG/KG	Endrin
4.1 U	UG/KG	Endosulfan II (beta)
4.1 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.1 U	UG/KG	Endosulfan Sulfate
4.1 Ü	UG/KG	4,4'-DDT (p,p'-DDT)
21 U	UG/KG	Methoxychlor
4.1 U	UG/KG	Endrin Ketone
4.1 U	UG/KG	Endrin Aldehyde
2.1 U	UG/KG	alpha-Chlordane /2
2.1 U	UG/KG	gamma-Chlordane /2
210 U	UG/KG	Toxaphene
41 U	UG/KG	PCB-1016 (Aroclor 1016)
84 U .	UG/KG	PCB-1221 (Aroclor 1221)
41 U	UG/KG	PCB-1232 (Aroclor 1232)
41 U	UG/KG	PCB-1242 (Aroclor 1242)
41 U	UG/KG	PCB-1248 (Aroclor 1248)
41 U	UG/KG	PCB-1254 (Aroclor 1254)
41 U	UG/KG	PCB-1260 (Aroclor 1260)
20	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Production Date: 06/20/2003 14:11

Sample 6765 FY 2003 Project: 03-0474

Produced by: Goddard, Denise Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 10:30

Pesticides Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS03SD /

Case No: 31635 MD No: 1XS5

Org Contractor: LIBRTY

D No: 1XS5 Media: SEDIMENT

Inorg Contractor: SENTIN

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
8.1 U	UG/KG	alpha-BHC
12	UG/KG	beta-BHC
8.1 U	UG/KG	delta-BHC
8.1 U	UG/KG	gamma-BHC (Lindane)
8.1 U	UG/KG	Heptachlor
8.1 U	UG/KG	Aldrin
8.1 U	UG/KG	Heptachlor Epoxide
8.1 U	UG/KG	Endosulfan I (alpha)
16 U	UG/KG	Dieldrin
16 U	UG/KG	4,4'-DDE (p,p'-DDE)
16 U	UG/KG	Endrin
16 U	UG/KG	Endosulfan II (beta)
16 U	UG/KG	4,4'-DDD (p,p'-DDD)
14 NJ_	UG/KG	Endosulfan Sulfate
16 U	UG/KG <sup>-</sup>	4,4'-DDT (p,p'-DDT)
81 U	UG/KG	Methoxychlor
16 U	UG/KG	Endrin Ketone
16 U	UG/KG	Endrin Aldehyde
8.1 U	UG/KG	alpha-Chlordane /2
8.1 U	UG/KG	gamma-Chlordane /2
810 U	UG/KG	Toxaphene
160 U	UG/KG	PCB-1016 (Aroclor 1016)
320 U	UG/KG	PCB-1221 (Aroclor 1221)
160 U	UG/KG	PCB-1232 (Aroclor 1232)
160 U	UG/KG	PCB-1242 (Aroclor 1242)
160 U	UG/KG	PCB-1248 (Aroclor 1248)
160 U	UG/KG	PCB-1254 (Aroclor 1254)
160 U	UG/KG	PCB-1260 (Aroclor 1260)
79	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Produced by: Goddard, Denise Sample 6766 FY 2003 Project: 03-0474 Requestor:

Pesticides Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS16SS /

MD No: 1XS6

Media: SURFACE SOIL (0" - 12")

Case No. 31635

**D No: 1XS6** 

Org Contractor: LIBRTY

Inorg Contractor: SENTIN

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

Project Leader: BSTRIGGO

Beginning: 04/22/2003 11:30

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2,2 U	UG/KG	Endosulfan I (alpha)
4.2 U	UG/KG	Dieldrin
2.5 J	UG/KG	4,4'-DDE (p,p'-DDE)
4.2 U	UG/KG	Endrin
4.2 U	UG/KG	Endosulfan II (beta)
4.2 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.2 U	UG/KG	Endosulfan Sulfate
4.2 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
14 J	UG/KG	Endrin Ketone
4.2 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U	UG/KG	Toxaphene
42 U	UG/KG	PCB-1016 (Aroclor 1016)
85 U	UG/KG	PCB-1221 (Aroclor 1221)
42 U	UG/KG	PCB-1232 (Aroclor 1232)
42 U	UG/KG	PCB-1242 (Aroclor 1242)
42 U	UG/KG	PCB-1248 (Aroclor 1248)
42 U	UG/KG	PCB-1254 (Aroclor 1254)
42 U	UG/KG	PCB-1260 (Aroclor 1260)
21	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

PESTIC	IDES/PCB	SAMPLE	ANAL	.YSIS

Production Date: 06/20/2003 14:11

Project: 03-0474 FY 2003 Sample 6767

Produced by: Goddard, Denise

**Pesticides Scan** 

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO

Beginning: 04/22/2003 11:40

Program: SF

Case No: 31635

Ending:

Id/Station: GS16SB /

MD No: 1XS7

Media: SUBSURFACE SOIL (> 12")

D No: 1XS7

Flowood, MS

inorg Contractor: SENTIN Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

		······································
RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
1.3 J	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
.2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.2 U	UG/KG	Dieldrin
4.2 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.2 U	UG/KG	Endrin
4:2 U	· UG/KG	Endosulfan II (beta)
4.2 U	· UG/KG	4,4'-DDD (p,p'-DDD)
4.2 U	UG/KG	Endosulfan Sulfate
4.2 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.2 U	UG/KG	Endrin Ketone
4.2 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U	UG/KG	Toxaphene
· 42 U	UG/KG	PCB-1016 (Aroclor 1016)
86 U	UG/KG	PCB-1221 (Aroclor 1221)
42 U	UG/KG	PCB-1232 (Aroclor 1232)
42 U	UG/KG	PCB-1242 (Aroclor 1242)
42 U	UG/KG	PCB-1248 (Aroclor 1248)
42 U	UG/KG	PCB-1254 (Aroclor 1254)
70	UG/KG	PCB-1260 (Aroclor 1260)
22	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

PESTICIDES/PCB SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:11

Project: 03-0474 FY 2003 Sample

Produced by: Goddard, Denise

Requestor:

Pesticides Scan

Project Leader: BSTRIGGO

Facility: Gulf States Creosoting

Flowood, MS

Beginning: 04/22/2003 12:05

Program: SF

Case No: 31635

Ending:

Id/Station: GS21SS / Media: SURFACE SOIL (0" - 12") MD No: 1XS8

Inorg Contractor: SENTIN

D No: 1XS8

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

		<del></del>
RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.2 U	UG/KG	Dieldrin
4.2 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.2 U	UG/KG	Endrin
4.2 U	UG/KG	Endosulfan II (beta)
4.2 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.2 U	UG/KG	Endosulfan Sulfate
4.2 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.2 U	UG/KG	Endrin Ketone
4.2 U	UG/KG	Endrin Aldehyde ,
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U -	UG/KG	Toxaphene
42 U	UG/KG	PCB-1016 (Aroclor 1016)
86 U	UG/KG	PCB-1221 (Aroclor 1221)
42 U	UG/KG	PCB-1232 (Aroclor 1232)
. 42 U	UG/KG	PCB-1242 (Aroclor 1242)
42 U	UG/KG	PCB-1248 (Aroclor 1248)
42 U	UG/KG	PCB-1254 (Aroclor 1254)
42 U	UG/KG	PCB-1260 (Aroclor 1260)
22	%	% Moisture
		!

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise Project: 03-0474 6769 FY 2003 Sample Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 12:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS21SB / MD No: 1XS9

Media: SUBSURFACE SOIL (> 12")		D No: 1XS9	Org Contractor: LIBRTY	
RESULTS 2.1 U	UNITS UG/KG	ANALYTE alpha-BHC		
2.1 U	UG/KG	beta-BHC		
2.1 U	UG/KG	delta-BHC		
2.1 U	UG/KG	gamma-BHC (Lindane)		
2.1 U	UG/KG	Heptachlor		
	· UG/KG	Aldrin		•
2.1 U	UG/KG	Heptachlor Epoxide		
2.1 U	UG/KG -	Endosulfan I (alpha)	•	
4.1 U	UG/KG	Dieldrin	•	
4.1 U	UG/KG	4,4'-DDE (p,p'-DDE)		
4.1 U	UG/KG	Endrin		
4.1 U	UG/KG	Endosulfan II (beta)		•
4.1 U	UG/KG	4,4'-DDD (p,p'-DDD)		
4.1 U	UG/KG	Endosulfan Sulfate		
4.1 U	UG/KG	4,4'-DDT (p,p'-DDT)	•	
21 U	UG/KG	Methoxychlor		
4.1 U	· UG/KG	Endrin Ketone	•	
4.1 U	UG/KG	Endrin Aldehyde		
2.1 U	UG/KG	alpha-Chlordane /2 gamma-Chlordane /2		
2.1 U 210 U	UG/KG UG/KG	Toxaphene		
41 U	UG/KG	PCB-1016 (Aroclor 1016)	1	
84 U	UG/KG	PCB-1221 (Aroclor 1221)		
41 U	UG/KG	PCB-1232 (Aroclor 1232)		
41 U	UG/KG	PCB-1242 (Aroclor 1242)		•
41 Ú	- UG/KG	PCB-1248 (Aroclor 1248)	•	
41 U	UG/KG	PCB-1254 (Aroclor 1254)	•	
41 U	UG/KG	PCB-1260 (Aroclor 1260)		
20	%	% Moisture	•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

6770

FY 2003

Project: 03-0474

Requestor:

Ending:

Produced by: Goddard, Denise

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:05

Program: SF

Case No: 31635

Id/Station; GS01GW / Media: GROUNDWATER

MD No: 1XT0 D No: 1XT0

Inorg Contractor: SENTIN

Org Contractor: A4

UNITS	ANALYTE
UG/L	alpha-BHC
UG/L	beta-BHC
UG/L	delta-BHC
UG/L	.gamma-BHC (Lindane);
UG/L	Heptachlor
UG/L	Aldrin
UG/L	Heptachlor Epoxide
UG/L	Endosulfan I (alpha)
UG/L	Dieldrin
UG/L	4,4'-DDE (p,p'-DDE)
UG/L	Endrin
	Endosulfan II (beta)
	4,4'-DDD (p,p'-DDD)
	Endosulfan Sulfate
UG/L	4,4'-DDT (p,p'-DDT)
	Methoxychlor
	Endrin Ketone
	Endrin Aldehyde
	alpha-Chlordane /2
	gamma-Chlordane /2
	Toxaphene
	PCB-1016 (Aroclor 1016)
	PCB-1221 (Aroclor 1221)
	PCB-1232 (Aroclor 1232)
	PCB-1242 (Aroclor 1242)
	PCB-1248 (Aroclor 1248)
	PCB-1254 (Aroclor 1254)
UG/L	PCB-1260 (Aroclor 1260)
	UG/L UG/L UG/L UG/L UG/L UG/L UG/L UG/L

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

PESTICIDES/PCB SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:11

Project: 03-0474 FY 2003 Sample 6771

Produced by: Goddard, Denise

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS01SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

MD No: 1XT1

D No: 1XT1

inorg Contractor: SENTIN

Org Contractor: LIBRTY

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.1 U	UG/KG	alpha-BHC
2.1 U	UG/KG	beta-BHC
2.1 U	UG/KG	delta-BHC
2.1 U	UG/KG	gamma-BHC (Lindane)
2.1 U	UG/KG	Heptachlor
2.1 Ú	UG/KG	Aldrin
2.1 U	UG/KG	Heptachlor Epoxide
2.1 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.0 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
4.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
21 U *	UG/KG	Methoxychlor
4.0 U	UG/KG	Endrin Ketone
4.0 U	UG/KG	Endrin Aldehyde
2.1 U	UG/KG	alpha-Chlordane /2
2.1 U	UG/KG	gamma-Chlordane /2
210 U	UG/KG	Toxaphene
40 U	UG/KG	PCB-1016 (Aroclor 1016)
82 U	UG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
40 U	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
18	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

PESTICIDES/PCB	SAMPLE	ANALYSIS

Production Date: 06/20/2003 14:11

DATA REPORTED ON DRY WEIGHT BASIS

Sample 6772 FY 2003 Project: 03-0474 Produced by: Goddard, Denise Requestor:
Pesticides Scan Project Leader: BSTRIGGO

Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:00
Program: SF Case No: 31635 Ending:

Id/Station: GS01SS / MD No: 1XT2 Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")		D No: 1XT2	Org Contractor: LIBRTY	
RESULTS	UNITS	ANALYTE		
2.0 U	UG/KG	alpha-BHC	•	
2.0 U	UG/KG	beta-BHC		
2.0 U	UG/KG	delta-BHC	•	
2.0 U	UG/KG	gamma-BHC (Lindane)		
2.0 U	UG/KG	Heptachlor		• •
2.0 U	UG/KG	Aldrin		
2.0 U -	UG/KG	Heptachlor Epoxide	•	•
2.0 U	UG/KG	Endosulfan I (alpha)		•
3.8 U	UĢ/KG	Dieldrin	•	
3.8 U	UG/KG	4,4'-DDE (p,p'-DDE)	•	
3.8 U	UG/KG	Endrin	•	•
3.8 U	UG/KG	Endosulfan II (beta)		
3.8 U	UG/KG	4,4'-DDD (p,p'-DDD)		
3.8 U	UG/KG	Endosulfan Sulfate	•	
3.8 U	UG/KG	4,4'-DDT (p,p'-DDT)		•
20 U	UG/KG	Methoxychlor		
3.8 U	UG/KG	Endrin Ketone		
	UG/KG	Endrin Aldehyde		
2.0 U	UG/KG	alpha-Chlordane /2		
2.0 U	UG/KG	gamma-Chlordane /2		•
200 U	UG/KG	Toxaphene		
38 U	UG/KG	PCB-1016 (Aroclor 1016	•	
78 U	UG/KG	PCB-1221 (Aroclor 1221)		
38 U	UG/KG	PCB-1232 (Aroclor 1232)	)	

PCB-1242 (Aroclor 1242)

PCB-1248 (Arocior 1248)

PCB-1254 (Aroclor 1254)

PCB-1260 (Aroclor 1260)

% Moisture

38 U

38 U

38 U

38 U

14

UG/KG

UG/KG

UG/KG

UG/KG

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample 6773 FY 2003 Project: 03-0474 Produced by: Goddard, Denise
Requestor:
Pesticides Scan
Facility: Gulf States Creosoting
Program: SF
Case No: 31635
Program: SF
Case No: 31635
Produced by: Goddard, Denise
Requestor:
Project Leader: BSTRIGGO
Beginning: 04/22/2003 13:15
Ending:

Id/Station: GS13SS / MD No: 1XT3 Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12") D No: 1XT3 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

· · · · · · · · · · · · · · · · · · ·					
RESULTS	UNITS	ANALYTE			
2.0 ປ	UG/KG	alpha-BHC			
2.0 U	UG/KG	beta-BHC			
2.0 ປ	UG/KG	delta-BHC			
2.0 U	UG/KG	gamma-BHC (Lindane)	•		
2.0 U	UG/KG	Heptachlor			•
2.0 U	UG/KG	Aldrin			
2.0 U	UG/KG	Heptachlor Epoxide			
2.0 U	UG/KG	Endosulfan I (alpha)			
3.8 U	UG/KG	Dieldrin			•
30 U	UG/KG	4,4'-DDE (p,p'-DDE)		•	
42 J	UG/KG	Endrin			
3.8 U	UG/KG	Endosulfan II (beta)	•		
3.8 U	UG/KG	4,4'-DDD (p,p'-DDD)			
3.8 U	UG/KG	Endosulfan Sulfate			
38 J	UG/KG	4,4'-DDT (p,p'-DDT)			
20 U	.UG/KG	Methoxychlor	•		
150	UG/KG	Endrin Ketone			
49 J	UG/KG	Endrin Aldehyde			
2.0 U	UG/KG	alpha-Chlordane /2			
18 U	UG/KG	gamma-Chlordane /2	•		
200 U	UG/KG	Toxaphene			
38 U	UG/KG	PCB-1016 (Aroclor 1016			
77 U	UG/KG	PCB-1221 (Aroclor 1221	)		
38 U	UG/KG	PCB-1232 (Aroclor 1232)			
38 U .	UG/KG	PCB-1242 (Aroclor 1242)	)		•
38 U	UG/KG	PCB-1248 (Aroclor 1248)			
38 U	UG/KG	PCB-1254 (Aroclor 1254			
38 U	UG/KG	PCB-1260 (Aroclor 1260)		-	
13	%	% Moisture			
		! !			

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/20/2003 14:11

Sample 6774 FY 2003 Project: 03-0474 Produced by: Goddard, Denise Requestor:

Pesticides Scan
Project Leader: BSTRIGGO
Facility: Gulf States Creosoting Flowood, MS
Beginning: 04/22/2003 13:25

Program: SF Case No: 31635 Ending: Id/Station: GS13SB / Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12") D No: 1XT4 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

RESULTS UNITS **ANALYTE** 2.0 U UG/KG alpha-BHC 2.0 U beta-BHC UG/KG delta-BHC 2.0 U UG/KG gamma-BHC (Lindane) 2.0 U UG/KG 2.0 U UG/KG Heptachlor 2.0 U UG/KG Aldrin Heptachlor Epoxide 2.0·U UG/KG 2.0 U UG/KG Endosulfan I (alpha) 3.9 U Dieldrin UG/KG 6.9 4,4'-DDE (p,p'-DDE) UG/KG 3.9 U UG/KG Endrin 3.9 U UG/KG Endosulfan II (beta) 3.9 U UG/KG 4,4'-DDD (p,p'-DDD) 3.9 U Endosulfan Sulfate UG/KG 3.9 U 4,4'-DDT (p,p'-DDT) UG/KG 20 U. UG/KG Methoxychlor 3.9 U UG/KG Endrin Ketone 3.9 U UG/KG Endrin Aldehyde 2.0 U alpha-Chlordane /2 UG/KG 2.0 U UG/KG gamma-Chlordane /2 200 U UG/KG Toxaphene 39 U UG/KG PCB-1016 (Aroclor 1016) 79 U UG/KG PCB-1221 (Aroclor 1221) 39 U UG/KG PCB-1232 (Aroclor 1232) 39 U UG/KG PCB-1242 (Aroclor 1242) 39 U UG/KG PCB-1248 (Aroclor 1248) 39 U PCB-1254 (Aroclor 1254) UG/KG 39 U UG/KG PCB-1260 (Aroclor 1260) 15 % Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Inora Contractor: SENTIN

Org Contractor: LIBRTY

Facility: Gulf States Creosoting

Program: SF

150 U

150 U

150 U

150 U

78

UG/KG

UG/KG

UG/KG

UG/KG

%

Production Date: 06/20/2003 14:11

Project: 03-0474 Sample 6775 FY 2003 Requestor: **Pesticides Scan** 

Case No: 31635

MD No: 1XT5

D No: 1XT5

Flowood, MS

Produced by: Goddard, Denise

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:35

Endina:

DATA REPORTED ON DRY WEIGHT BASIS

Id/Station: GS04SD / Media: SEDIMENT ANALYTE **RESULTS UNITS** alpha-BHC 7.7 U UG/KG beta-BHC .7.7 U UG/KG 7.7 U UG/KG delta-BHC 7.7 U gamma-BHC (Lindane) UG/KG Heptachlor 7.7 U UG/KG 7.7 U UG/KG Aldrin Heptachlor Epoxide 7.7 U· UG/KG 7.7 U Endosulfan I (alpha) UG/KG 15 U Dieldrin UG/KG 15 U UG/KG 4,4'-DDE (p,p'-DDE) 15 U UG/KG Endrin 15 U UG/KG Endosulfan II (beta) 15 U UG/KG 4,4'-DDD (p,p'-DDD) Endosulfan Sulfate 15 U UG/KG 15 U UG/KG 4,4'-DDT (p,p'-DDT) 77 U Methoxychlor UG/KG Endrin Ketone 15 U UG/KG Endrin Aldehyde 15 U UG/KG 7.7 U UG/KG alpha-Chlordane /2 gamma-Chlordane /2 7.7 U UG/KG 770 U UG/KG Toxaphene 150 U UG/KG PCB-1016 (Aroclor 1016) PCB-1221 (Aroclor 1221) 300 U UG/KG PCB-1232 (Aroclor 1232) 150 U UG/KG

PCB-1242 (Aroclor 1242)

PCB-1248 (Aroclor 1248)

PCB-1254 (Aroclor 1254)

PCB-1260 (Aroclor 1260)

% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value;

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

6776 FY 2003 Project: 03-0474 Sample

Pesticides Scan

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS05SD / Media: SEDIMENT

Flowood, MS

Case No: 31635

MD No: 1XT6 D No: 1XT6

Org Contractor: LIBRTY

Inorg Contractor: SENTIN

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:15

Endina:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
3.7 U	UG/KG	alpha-BHC
3.7 U	UG/KG	beta-BHC
3.7 U	UG/KG	delta-BHC
3.7 U	UG/KG	gamma-BHC (Lindane)
3.7 U	UG/KG	Heptachlor
3.7 U	UG/KG	Aldrin
3.7 U	UG/KG	Heptachlor Epoxide
3.7 U	UG/KG	Endosulfan I (alpha)
7.2 U	UG/KG	Dieldrin
7.2 U	UG/KG `	4,4'-DDE (p,p'-DDE)
7.2 U	UG/KG	Endrin
7.2 U	UG/KG	Endosulfan II (beta)
7.2 U	UG/KG	4,4'-DDD (p,p'-DDD)
7.2 U	UG/KG	Endosulfan Sulfate
7.2 U	UG/KG	4,4'-DDT (p,p'-DDT)
37 U	UG/KG	Methoxychlor
7.2 U	UG/KG	Endrin Ketone
7.2 U	UG/KG	Endrin Aldehyde
3.7 U	UG/KG	alpha-Chlordane /2
3.7 U	UG/KG	gamma-Chlordane /2
370 U	UG/KG	Toxaphene
72 U	UG/KG	PCB-1016 (Aroclor 1016)
150 U	UG/KG	PCB-1221 (Aroclor 1221)
72 U	UG/KG	PCB-1232 (Aroclor 1232)
72 U .	UG/KG	PCB-1242 (Aroclor 1242)
72 U	UG/KG	PCB-1248 (Aroclor 1248)
72 U	UG/KG	PCB-1254 (Aroclor 1254)
72 U	UG/KG	PCB-1260 (Aroclor 1260)
54	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample **Pesticides Scan** 

6777

Facility: Gulf States Creosoting

FY 2003 Project: 03-0474 Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 12:45

Ending:

Program: SF

Flowood, MS

Case No: 31635

Id/Station: GS07SD / Media: SEDIMENT

MD No: 1XT7

Inorg Contractor: SENTIN D No: 1XT7

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.1 U	UG/KG	alpha-BHC
2.1 U	UG/KG	beta-BHC
2.1 U	UG/KG	delta-BHC
2.1 U	UG/KG	gamma-BHC (Lindane)
2.1 U	UG/KG	Heptachlor
2.1. U	UG/KG	Aldrin
2.1 U	UG/KG	Heptachlor Epoxide
2.1 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.0 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
4.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
21 U .	UG/KG	Methoxychlor
4.0 U	UG/KG	Endrin Ketone
4.0 Ü	UG/KG	Endrin Aldehyde
2.1 U	UG/KG	alpha-Chlordane /2
2.1 U	UG/KG	gamma-Chlordane /2
210 U	UG/KG	Toxaphene
40 U	UG/KG	PCB-1016 (Aroclor 1016)
82 U	UG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
40 U	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
18	%	% Moisture
		ļ

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample 6778 FY 2003 Project: 03-0474

**Pesticides Scan** 

Facility: Gulf States Creosoting

Program: SF Id/Station: GS23SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635 MD No: 1XT8

D No: 1XT8

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:55

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

			·	
RESULTS	UNITS	ANALYTE		:
1.7 NJ	UG/KG	alpha-BHC		
2.3 U	UG/KG	beta-BHC		
2.3 U	UG/KG	delta-BHC	·	
2.3 U	UG/KG	gamma-BHC (Lindane)		
2.3 U	UG/KG	Heptachlor		
2.3 U	UG/KG	Aldrin	•	
2.3 U	UG/KG	Heptachlor Epoxide	1	
2.3 U	UG/KG	Endosulfan I (alpha)		
4.5 U	UG/KG	Dieldrin	! 	
4.5 U	UG/KG	4,4'-DDE (p,p'-DDE)	1	
4.5 U	UG/KG	Endrin	1	
4.5 U	UG/KG	Endosulfan II (beta)	!	
4.5 U	UG/KG	4,4'-DDD (p,p'-DDD)	·	
4.5 U	UG/KG	Endosulfan Sulfate	İ	
4.5 U	UG/KG	4,4'-DDT (p,p'-DDT)		
23 U	UG/KG	Methoxychlor		
6.9·J	UG/KG	Endrin Ketone		
4.5 U	UG/KG	Endrin Aldehyde		
2.3 U	UG/KG	alpha-Chlordane /2		
2.3 U	UG/KG ·	gamma-Chlordane /2		
230 U	UG/KG	Toxaphene		
45 U	UG/KG	PCB-1016 (Aroclor 101	<b>6</b> )	
91 U	UG/KG	PCB-1221 (Aroclor 122	1)	
45 U	UG/KG	PCB-1232 (Aroclor 123)	<b>2</b> )	
45 U	UG/KG	PCB-1242 (Aroclor 124)	2)	
45 U	UG/KG	PCB-1248 (Aroclor 124)	8)	
45 U	UG/KG	PCB-1254 (Aroclor 1254)	4)	
45 U	UG/KG	PCB-1260 (Aroclor 126)	Ó)	
26	%	% Moisture		
•				

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

DECTION DEC/DA	OD CAMEDIE	ANIAL VOIC
PESTICIDES/PO	'R SAMLLE	AIVALTSIS

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise FY 2003 Sample Project: 03-0474 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:40 Program: SF Case No: 31635 Ending: . Id/Station: GS15SS / Inorg Contractor: SENTIN MD No: 1XT9 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT9 DATA REPORTED ON DRY WEIGHT BASIS

	RESULTS	UNITS	ANALYTE	
	2.0 U	UG/KG	alpha-BHC	
	2.0 U	UG/KG	beta-BHC	
	2.0 U	UG/KG	delta-BHC	
	2.0 U	UG/KG	gamma-BHC (Lindane)	
	2.0 U	UG/KG	Heptachlor	
	2.0 U	UG/KG	Aldrin	
	2.0 U	UG/KG	Heptachlor Epoxide	
	2.0 U	UG/KG	Endosulfan I (alpha)	
٠	3.9 U	UG/KG	Dieldrin	
	- 3.9 U	UG/KG	4,4'-DDE (p,p'-DDE)	
	11 U	UG/KG	Endrin	
	3.9 U	UG/KG	Endosulfan II (beta)	
	3.9 U	UG/KG	4,4'-DDD (p,p'-DDD)	
	3.9 U	UG/KG	Endosulfan Sulfate	
	30 NJ	UG/KG	4,4'-DDT (p,p'-DDT)	
	61 U	UG/KG	Methoxychlor	
	3.9 U	UG/KG	Endrin Ketone	
	. 27 NJ	UG/KG	Endrin Aldehyde	
	2.0 U	UG/KG	alpha-Chlordane /2	
	2.0 U	UG/KG	gamma-Chlordane /2	
	200 U	UG/KG	Toxaphene	
	39 U	UG/KG	PCB-1016 (Aroclor 1016	5)
	80 U	UG/KG	PCB-1221 (Aroclor 1221	
	39 U	UG/KG	PCB-1232 (Aroclor 1232	<u>,</u>
	39 U	UG/KG	PCB-1242 (Aroclor 1242	
	39 U 🔭	UG/KG	PCB-1248 (Aroclor 1248	<b>3)</b>
	39 U	UG/KG	PCB-1254 (Arocior 1254	
	39 U	UG/KG	PCB-1260 (Aroclor 1260	))
	16	%	% Moisture	
			•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

PESTICIDES/PCB SAMPLE ANAL'	YSIS
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Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise Project: 03-0474 Sample 6780 FY 2003 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood MS Beginning: 04/22/2003 13:50 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS15SB / MD No: 1XW0 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XW0 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS** ANALYTE 2.2 U UG/KG alpha-BHC 2.2 U UG/KG beta-BHC 2.2 U delta-BHC UG/KG 2.2 U UG/KG gamma-BHC (Lindane) 2.2 U UG/KG Heptachlor 2.2 U UG/KG Aldrin 2.2 U UG/KG Heptachlor Epoxide 2.2 U UG/KG Endosulfan I (alpha) 4.2 U UG/KG Dieldrin 4.2 U UG/KG 4,4'-DDE (p,p'-DDE) 4.2 U UG/KG Endrin 4.2 U UG/KG Endosulfan II (beta) 4.2 U UG/KG 4,4'-DDD (p,p'-DDD) 4.2 U UG/KG Endosulfan Sulfate 4.2 U 4,4'-DDT (p,p'-DDT) UG/KG 22 U UG/KG Methoxychlor 4.2 U ° UG/KG **Endrin Ketone** 4.2 U Endrin Aldehyde UG/KG 2.2 U UG/KG alpha-Chlordane /2 2.2 U UG/KG gamma-Chlordane /2 220 U UG/KG Toxaphene 42 U UG/KG PCB-1016 (Aroclor 1016) PCB-1221 (Aroclor 1221) 85 U UG/KG 42 U UG/KG PCB-1232 (Aroclor 1232) PCB-1242 (Aroclor 1242) 42 U UG/KG PCB-1248 (Aroclor 1248) 42 U UG/KG 42 U UG/KG PCB-1254 (Aroclor 1254) 42 U UG/KG PCB-1260 (Aroclor 1260) 21 % Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

4.3 U

4.3 U

4.3 U

4.3 U

4.3 U

4.3 U

22 U

4.3 U

4.3 U

2.2 U

2.2 U

220 U

43 U

87 U

43 U

43 U

43 U

43 U

43 U

23

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

UG/KG

%

4.4'-DDE (p.p'-DDE)

Endosulfan II (beta)

4,4'-DDD (p,p'-DDD)

Endosulfan Sulfate

4,4'-DDT (p,p'-DDT)

alpha-Chlordane /2

gamma-Chlordane /2

PCB-1016 (Aroclor 1016) PCB-1221 (Aroclor 1221)

PCB-1232 (Aroclor 1232) PCB-1242 (Aroclor 1242)

PCB-1248 (Aroclor 1248)

PCB-1254 (Aroclor 1254)

PCB-1260 (Aroclor 1260)

Methoxychlor

**Endrin Ketone** 

Toxaphène

% Moisture

Endrin Aldehyde

Endrin

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise Project: 03-0474 Sample 6781 FY 2003 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:13 Program: SF Case No: 31635 Endina: Id/Station: GS02SS / MD No: 1XW1 Inorg Contractor: SENTIN Org Contractor: LIBRTY D No: 1XW1 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** 2.2 U UG/KG alpha-BHC 2.2 U UG/KG beta-BHC 2.2 U delta-BHC UG/KG 2.2 U UG/KG gamma-BHC (Lindane) 2.2 U UG/KG Heptachlor 2.2 U UG/KG Aldrin 2.2 U UG/KG Heptachlor Epoxide 2.2 U -UG/KG Endosulfan I (alpha) 4.3 U UG/KG Dieldrin

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable

PESTICIDES/PCB	SAMPLE.	ANALYSIS

Production Date: 06/20/2003 14:11

Sample 6782 FY 2003 Project: 03-0474 Produced by: Goddard, Denise
Requestor:
Pesticides Scan
Facility: Gulf States Creosoting Flowood, MS Project Leader: BSTRIGGO
Beginning: 04/22/2003 13:34

Program: SF Case No: 31635 Ending: Id/Station: GS02SB / Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12") D No: 1XW2 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

			<del></del>	
RESULTS	UNITS	ANALYTE		
2.1 U	UG/KG	alpha-BHC		
2.1 U	UG/KG	beta-BHC		
2.1 U	UG/KG	delta-BHC		
2.1 U	UG/KG	gamma-BHC (Lindane)		
2.1 U	UG/KG	Heptachlor		
2.1 U	UG/KG	Aldrin		
2.1 U	UG/KG	Heptachlor Epoxide		
2.1 U	UG/KG	Endosulfan I (alpha)		
4.1 U	UG/KG	Dieldrin		
- 4.1 U	UG/KG	4,4'-DDE (p,p'-DDE)		
4.1 U	UG/KG	Endrin		
4.1 U	UG/KG	Endosulfan II (beta)		
4.1 U	UG/KG	4,4'-DDD (p,p'-DDD)		
4.1 U	UG/KG	Endosulfan Sulfate		
. 4.1 U	UG/KG	4,4'-DDT (p,p'-DDT)		•
21 U	UG/KG	Methoxychlor		
4.1 U	UG/KG	Endrin Ketone		
4.1 U	UG/KG	Endrin Aldehyde		
2.1 U	UG/KG	alpha-Chlordane /2		
2.1 U	UG/KG	gamma-Chlordane /2		
210 U	UG/KG	Toxaphene		
41 U	UG/KG	PCB-1016 (Aroclor 101		
83 U	UG/KG	PCB-1221 (Aroclor 122		
41 U	UG/KG	PCB-1232 (Aroclor 123)		
41 U	UG/KG	PCB-1242 (Aroclor 124)		
41 U 、		PCB-1248 (Aroclor 124)		
41 U	UG/KG	PCB-1254 (Aroclor 125		
41 U	UG/KG	PCB-1260 (Aroclor 126	<i>(</i> )	
19	%	% Moisture		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

PESTICIDES/PCB SAMPLE ANALYSIS	
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Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise Project: 03-0474 Sample 6783 FY 2003 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 14:40 Program: SF Case No: 31635 Ending: Id/Station: GS14SS / MD No: 1XW3 Inorg Contractor: SENTIN Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XW3 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
2.0 U	UG/KG	Heptachior
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
3.8 U	UG/KG	Dieldrin
3.8 U	UG/KG	4,4'-DDE (p,p'-DDE)
3.8 U	UG/KG	Endrin
3.8 U	UG/KG	Endosulfan II (beta)
3.8 U	UG/KG	4,4'-DDD (p,p'-DDD)
3.8 U	UG/KG	Endosulfan Sulfate
10 NJ	UG/KG	4,4'-DDT (p,p'-DDT)
34 NJ	UG/KG	Methoxychlor
3.8 U	UG/KG	Endrin Ketone
3.8 U	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
38 U	UG/KG	PCB-1016 (Aroclor 1016)
78 U	UG/KG	PCB-1221 (Aroclor 1221)
38 U	UG/KG	PCB-1232 (Aroclor 1232)
38 U	UG/KG	PCB-1242 (Aroclor 1242)
38 U	UG/KG	PCB-1248 (Aroclor 1248)
38 U	UG/KG	PCB-1254 (Aroclor 1254)
38 U	UG/KG	PCB-1260 (Aroclor 1260)
14	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

<b>PESTICIDES</b>	/PCR	SAMPL	FAN	AL YSIS
PESTICIDES	/rub	SHIVIFL	$-\sim$	AL I JIJ

Production Date: 06/20/2003 14:11

Sample 6784 FY 2003 Project: 03-0474

Produced by: Goddard, Denise

**Pesticides Scan** 

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:50

Program: SF

Case No: 31635

ld/Station: GS14SB /

MD No: 1XW4 Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

Inorg Contractor: SENTIN
Org Contractor: LIBRTY

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
2.0 U	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
3.9 U	UG/KG	Dieldrin
3.9 U	UG/KG	4,4'-DDE (p,p'-DDE)
3.9 U	UG/KG	Endrin
3.9 U	UG/KG	Endosulfan II (beta)
3.9 U	UG/KG	4,4'-DDD (p,p'-DDD)
3.9 U	UG/KG	Endosulfan Sulfate
16 J	UG/KG	4,4'-DDT (p,p'-DDT)
43 U	UG/KG	Methoxychlor
3.9 U	UG/KG	Endrin Ketone
3.9 U	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
39 U	UG/KG	PCB-1016 (Aroclor 1016)
79 U	UG/KG	PCB-1221 (Aroclor 1221)
39 U	UG/KG	PCB-1232 (Aroclor 1232)
39 U	UG/KG	PCB-1242 (Aroclor 1242)
39 U	UG/KG	PCB-1248 (Aroclor 1248)
39 U	UG/KG	PCB-1254 (Aroclor 1254)
39 U	UG/KG	PCB-1260 (Aroclor 1260)
15	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

Flowood, MS

D No: 1XW4

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

### PESTICIDES/PCB SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:11

Project: 03-0474 Sample 6785 FY 2003

Produced by: Goddard, Denise

**Pesticides Scan** 

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO

Program: SF

Beginning: 04/22/2003 15:05

Case No: 31635

Id/Station: GS10SS /

Inorg Contractor: SENTIN

Media: SURFACE SOIL (0" - 12")

MD No: 1XW5 Org Contractor: LIBRTY D No: 1XW5

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
1.9 U	UG/KG	alpha-BHC
12 U	UG/KG	beta-BHC
1.9 U	UG/KG	delta-BHC
1.9 U	UG/KG	gamma-BHC (Lindane)
1.9 U	UG/KG	Heptachlor
1.9 U	UG/KG	Aldrin
1.9 U	UG/KG	Heptachlor Epoxide
1.9 U	UG/KG	Endosulfan I (alpha)
3.8 U	UG/KG	Dieldrin
3.8 U	UG/KG	4,4'-DDE (p,p'-DDE)
15 U	UG/KG	Endrin
3.8 U	UG/KG	Endosulfan II (beta)
3.8 U	UG/KG	4,4'-DDD (p,p'-DDD)
3.8 U	UG/KG	Endosulfan Sulfate
43 J	UG/KG	4,4'-DDT (p,p'-DDT)
81 NJ	UG/KG	Methoxychlor
3.8 U	UG/KG	Endrin Ketone
3.8 U	UG/KG	Endrin Aldehyde
1.9 U	UG/KG	alpha-Chlordane /2
1.9 U	UG/KG	gamma-Chlordane /2
190 U	UG/KG	Toxaphene
38 U	UG/KG	PCB-1016 (Aroclor 1016)
76 U	UG/KG	PCB-1221 (Aroclor 1221)
38 U	UG/KG	PCB-1232 (Aroclor 1232)
38 U	UG/KG	PCB-1242 (Aroclor 1242)
38 U	UG/KG	PCB-1248 (Aroclor 1248)
38 U -	UG/KG	PCB-1254 (Aroclor 1254)
38 U	UG/KG	PCB-1260 (Aroclor 1260)
12	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

Flowood, MS

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

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DEATIONEO/DAD	CARROLE ANIAL VOIC
PESTICIDES/PCB	SAMPLE ANALYSIS

Ending:

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample : 6786 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 15:15 Program: SF Case No: 31635

Inorg Contractor: SENTIN Id/Station: GS10SB / MD No: 1XW6 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No. 1XW6 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
19 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
2.0 U	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
37 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
8.7 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
110 N	UG/KG	4,4'-DDT (p,p'-DDT)
360 U	UG/KG	Methoxychlor
180 N	UG/KG	Endrin Ketone
15 U .	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
40 U	UG/KG	PCB-1016 (Aroclor 1016)
81 U	UG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
40 U	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
17	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

6787

FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:55

Ending:

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

MD No: 1XW7

Inorg Contractor: SENTIN

Id/Station: GS02GW / Media: GROUNDWATER

D No: 1XW7

Org Contractor: A4

RESULTS	UNITS	ANALYTE
0.010 U	UG/L	alpha-BHC
0.010 U	UG/L	beta-BHC
0.010 U	UG/L	delta-BHC
0.010 U	UG/L	gamma-BHC (Lindane)
0.010 U	UG/L	Heptachlor
0.010 U	UG/L	Aldrin
0.010 U	UG/L	Heptachior Epoxide
0.010 U	UG/L	Endosulfan I (alpha)
0.020 U ·	UG/L	Dieldrin
0.020 UJ	UG/L	4,4'-DDE (p,p'-DDE)
0.020 UJ	UG/L	Endrin
0.020 U	UG/L	Endosulfan II (beta)
0.020 UJ	UG/L	4,4'-DDD (p,p'-DDD)
0.020 U	UG/L	Endosulfan Sulfate
0.020 UJ	UG/L	4,4'-DDT (p,p'-DDT)
0.10 U	UG/L	Methoxychlor
0.020 UJ	UG/L	Endrin Ketone
0.020 UJ	UG/L	Endrin Aldehyde
0.010 U	UG/L	alpha-Chlordane /2
0.010 U	UG/L	gamma-Chlordane /2
1.0 U	UG/L	Toxaphene
0.20 U	UG/L	PCB-1016 (Aroclor 1016)
0.40 U	UG/L	PCB-1221 (Aroclor 1221)
0.20 U	UG/L	PCB-1232 (Aroclor 1232)
0.20 U	UG/L	PCB-1242 (Aroclor 1242)
0.20 U	UG/L	PCB-1248 (Aroclor 1248)
0.20 U	UG/L	PCB-1254 (Aroclor 1254)
0.20 U	UG/L	PCB-1260 (Aroclor 1260)
		!

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474 6788 FY 2003 Sample

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS11SS /

Media: SURFACE SOIL (0" - 12")

Case No: 31635

MD No: 1XW8 D No: 1XW8

Inorg Contractor: SENTIN

**Org Contractor: LIBRTY** 

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:00

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
1.8 U	UG/KG	alpha-BHC
5.9 U	UG/KG	beta-BHC
1.8 U	UG/KG	delta-BHC
1.8 U	UG/KG	gamma-BHC (Lindane)
1.8 U	UG/KG	Heptachlor
1.8 U	UG/KG	Aldrin
1.8 U	UG/KG	Heptachlor Epoxide
1.8 U	UG/KG	Endosulfan I (alpha)
3.5 U	UG/KG	Dieldrin
3.5 U	UG/KG	4,4'-DDE (p,p'-DDE)
3.5 U	UG/KG	Endrin
3.5 U	UG/KG	Endosulfan II (beta)
· 3.5 U	UG/KG	4,4'-DDD (p,p'-DDD)
3.5 U	UG/KG	Endosulfan Sulfate
30 J	UG/KG	4,4'-DDT (p,p'-DDT)
54 U	UG/KG	Methoxychlor
3.5 U	UG/KG	Endrin Ketone
3.5 U	UG/KG	Endrin Aldehyde
1.8 U	UG/KG	alpha-Chlordane /2
1.8 U	UG/KG	gamma-Chlordane /2
180 U	UG/KG	Toxaphene
35 U	UG/KG	PCB-1016 (Aroclor 1016)
72 U	UG/KG	PCB-1221 (Aroclor 1221)
35 U	UG/KG	PCB-1232 (Aroclor 1232)
35 U	UG/KG	PCB-1242 (Aroclor 1242)
35 U	UG/KG	PCB-1248 (Aroclor 1248)
35 U	UG/KG	PCB-1254 (Aroclor 1254)
35 U *	UG/KG	PCB-1260 (Aroclor 1260)
7.0	%	% Moisture
		!
		1

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value...

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474 Sample . 6789 FY 2003

**Pesticides Scan** 

Facility: Gulf States Creosoting

Program: SF

ld/Station: GS11SB /

Media: SUBSURFACE SOIL (> 12")

Flowood, MS

Case No: 31635 MD No: 1XW9

D No: 1XW9

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:10

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
2.0 U	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.0 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
4.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
20 U	UG/KG	Methoxychlor
4.0 U	UG/KG	Endrin Ketone
4.0 U	UG/KG	Endrin Aldehyde
2.0 U -	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
40 U	UG/KG	PCB-1016 (Aroclor 1016)
81 U	UG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
40 U	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
17	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Produced by: Goddard, Denise Project: 03-0474 Sample 6790 FY 2003 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:15 Program: SF Case No: 31635 Ending: Id/Station: GS09SS / Inorg Contractor: SENTIN MD No: 1XX0 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX0 DATA REPORTED ON DRY WEIGHT BASIS

 			·	
RESULTS	UNITS	ANALYTE		•
2.3 U	UG/KG	alpha-BHC		
2.3 U	UG/KG	beta-BHC	•	
2.3 U	UG/KG	delta-BHC		
2.3 U	UG/KG	gamma-BHC (Lindane)		
2.3 U	UG/KG	Heptachlor		
2.3 U	UG/KG	Aldrin		
2.3 U	UG/KG	Heptachlor Epoxide		
2.3 U	UG/KG	Endosulfan I (alpha)		
4.5 U .	UG/KG	Dieldrin		•
4.5 U	UG/KG	4,4'-DDE (p,p'-DDE)		
4.5 U	UG/KG	Endrin		
4.5 U	UG/KG	Endosulfan II (beta)		
4.5 U	UG/KG	4,4'-DDD (p,p'-DDD)		
4.5 U	UG/KG	Endosulfan Sulfate		
4.5 U	UG/KG	4,4'-DDT (p,p'-DDT)		
23 U	UG/KG	Methoxychlor	·	
4.5 U	UG/KG	Endrin Ketone		
4.5 U	UG/KG	Endrin Aldehyde	1	
2.3 U	UG/KG	alpha-Chlordane /2		. •
2.3 U	UG/KG	gamma-Chlordane /2		
230 U	UG/KG	Toxaphene	i I	
45 U	UG/KG	PCB-1016 (Aroclor 1016		
91 U	UG/KG	PCB-1221 (Aroclor 1221		•
45 U	UG/KG	PCB-1232 (Aroclor 1232		
45 U	UG/KG	PCB-1242 (Aroclor 1242		
45 U	UG/KG	PCB-1248 (Aroclor 1248		
45 U	UG/KG	PCB-1254 (Aroclor 1254		
45 U	UG/KG	PCB-1260 (Aroclor 1260	<b>)</b> )	
26	%	% Moisture		
		•		

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

FY 2003 Project: 03-0474 Sample

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635 Id/Station: GS09SB /

Media: SUBSURFACE SOIL (> 12")

MD No: 1XX1

Inorg Contractor: SENTIN D No: 1XX1

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:25

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.2 U	UG/KG	Dieldrin
4.2 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.2 U	UG/KG	Endrin
4.2 U	UG/KG	Endosulfan II (beta)
4.2 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.2 U	UG/KG	Endosulfan Sulfate
4.2 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.2 U	UG/KG	Endrin Ketone
4.2 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U	UG/KG	Toxaphene
42 U	UG/KG	PCB-1016 (Aroclor 1016)
85 U	UG/KG	PCB-1221 (Aroclor 1221)
42 U	UG/KG	PCB-1232 (Aroclor 1232)
42 U	UG/KG	PCB-1242 (Aroclor 1242)
42 U	UG/KG	PCB-1248 (Aroclor 1248)
42 U	UG/KG	PCB-1254 (Aroclor 1254)
42 U 🕠	UG/KG	PCB-1260 (Aroclor 1260)
21 <sup>.</sup>	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

PESTICIDES/PCB SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:11

FY 2003 Project: 03-0474 Sample

Produced by: Goddard, Denise

Pesticides Scan

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO

Program: SF

Case No: 31635

Beginning: 04/22/2003 15:15

Id/Station: GS06SD /

Flowood, MS

Endina:

Media: SEDIMENT

MD No: 1XX2 Inorg Contractor: SENTIN Org Contractor: LIBRTY D No: 1XX2

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.1 U	UG/KG	alpha-BHC
2.1 U	UG/KG	beta-BHC
2.1 U	UG/KG	delta-BHC
2.1 U	UG/KG	gamma-BHC (Lindane)
2.1 U	UG/KG	Heptachlor
2.1 U	UG/KG	Aldrin
2.1 U	UG/KG	Heptachlor Epoxide
2.1 U	UG/KG	Endosulfan I (alpha)
4.1 U	UG/KG	Dieldrin
4.1 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.1 U	UG/KG	Endrin
4.1 U	UG/KG	Endosulfan II (beta)
4.1 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.1 U .	UG/KG	Endosulfan Sulfate
4.1 U	UG/KG	4,4'-DDT (p,p'-DDT)
21 U	UG/KG	Methoxychlor
4.1 U	UG/KG	Endrin Ketone
4.1 U	UG/KG .	Endrin Aldehyde
2.1 U .	UG/KG	alpha-Chlordane /2
2.1 U	UG/KG	gamma-Chlordane /2
210 U	UG/KG	Toxaphene
41 U	UG/KG	PCB-1016 (Aroclor 1016)
84 U	UG/KG	PCB-1221 (Aroclor 1221)
41 U	UG/KG	PCB-1232 (Aroclor 1232)
41 U	UG/KG	PCB-1242 (Aroclor 1242)
41 U	UG/KG	PCB-1248 (Aroclor 1248)
41 U	UG/KG	PCB-1254 (Aroclor 1254)
41 U	UG/KG	PCB-1260 (Aroclor 1260)
20	%	% Moisture
		1

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. | an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample FY 2003 Project: 03-0474

Pesticides Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

ld/Station: GS08SD / Media: SEDIMENT

MD No: 1XX3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY D No: 1XX3

Ending:

Requestor:

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise

Project Leader: BSTRIGGO

Beginning: 04/22/2003 15:50

RESULTS	UNITS	ANALYTE
2.1 U	UG/KG	alpha-BHC
2.1 U	UG/KG	beta-BHC
2.1 U	UG/KG	delta-BHC
2.1 U	UG/KG	gamma-BHC (Lindane)
2.1 U	UG/KG	Heptachlor
2.1 U	UG/KG	Aldrin
2.1 U	UG/KG	Heptachlor Epoxide
2.1 U	UG/KG	Endosulfan I (alpha)
4.1 U	UG/KG	Dieldrin
4.1 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.1 U	UG/KG	Endrin
4.1 U	UG/KG	Endosulfan II (beta)
4.1 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.1 U	UG/KG	Endosulfan Sulfate
4.1 U	UG/KG	4,4'-DDT (p,p'-DDT)
21 U	UG/KG	Methoxychlor
4.1 Ü	UG/KG	Endrin Ketone
4.1 U	UG/KG	Endrin Aldehyde
2.1 U	UG/KG	alpha-Chlordane /2
2.1 U	UG/KG	gamma-Chlordane /2
210 U	UG/KG	Toxaphene
41 U	UG/KG	PCB-1016 (Aroclor 1016)
83 U	UG/KG	PCB-1221 (Aroclor 1221)
41 U	UG/KG	PCB-1232 (Aroclor 1232)
41 U	UG/KG	PCB-1242 (Aroclor 1242)
41 U.	UG/KG	PCB-1248 (Aroclor 1248)
41 U	UG/KG	PCB-1254 (Aroclor 1254)
41 U	UG/KG	PCB-1260 (Aroclor 1260)
19	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

PESTICIDES/PCB	CAMDIE	ANAL VOIC
PESTICIDES/PCD	SAMPLE	ANALTSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise Sample 6794 FY 2003 Project: 03-0474 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:42 Program: SF Case No: 31635 Ending: Id/Station: GS22SS / MD No: 1XX4 Inorg Contractor: SENTIN Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX4 DATA REPORTED ON DRY WEIGHT BASIS

Modia.	" NOL 00	/IL (0 12)	D 140: 17004
RESULTS	UNITS	ANALYTE	·
2.0 U-	UG/KG	alpha-BHC	•
2.0 U	UG/KG	beta-BHC	
2.0 U	UG/KG	delta-BHC	•
2.0 U	UG/KG	gamma-BHC (Lindane)	
2.0 U	UG/KG	Heptachlor	
2.0 U	UG/KG	Aldrin	
2.0 U	UG/KG	Heptachlor Epoxide	
2.0 U	UG/KG	Endosulfan I (alpha)	,
3.8 U	UG/KG	Dieldrin	
3.8 U	UG/KG	4,4'-DDE (p,p'-DDE)	
3.8 U	UG/KG	Endrin	-
3.8 U	UG/KG	Endosulfan II (beta)	:
3.8 U	UG/KG	4,4'-DDD (p,p'-DDD)	
3.8 U	UG/KG	Endosulfan Sulfate	
3.8 U	UG/KG	4,4'-DDT (p,p'-DDT)	
20 U	UG/KG	Methoxychlor	
3.8 U	UG/KG	Endrin Ketone	
3.8 U	UG/KG	Endrin Aldehyde	
2.0 U	UG/KG	alpha-Chlordane /2	
2.0 U	UG/KG	gamma-Chlordane /2	! !
200 U	UG/KG	Toxaphene	} <b>-</b>
38 U	UG/KG	PCB-1016 (Aroclor 101)	
. 77 U	UG/KG	PCB-1221 (Aroclor 122)	
38 U	UG/KG	PCB-1232 (Aroclor 1232	
38 U	UG/KG	PCB-1242 (Aroclor 1242	•
38 U	UG/KG	PCB-1248 (Aroclor 1248	
38 U	UG/KG	PCB-1254 (Aroclor 1254	
38 U	UG/KG	PCB-1260 (Aroclor 1260	<u>!)</u>
13	%	% Moisture	i

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Production Date: 06/20/2003 14:11

Project: 03-0474 Sample 6795 FY 2003

**Pesticides Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS06SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635

MD No: 1XX5

D·No: 1XX5

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:40

Endina:

DATA REPORTED ON DRY WEIGHT BASIS

F	RESULTS	UNITS	ANALYTE	
	2.3 U	UG/KG	alpha-BHC	
	2.3 U	UG/KG	beta-BHC	
	2.3 U	UG/KG	delta-BHC	
	2.3 U	UG/KG	gamma-BHC (Lindane)	
	2.3 U	UG/KG	Heptachlor	
	2.3 U	UG/KG	Aldrin	
	2.3 U	UG/KG	Heptachlor Epoxide	
	2.3 U	UG/KG	Endosulfan I (alpha)	
	4.5 U	UG/KG	Dieldrin	
	4.5 U	UG/KG	4,4'-DDE (p,p'-DDE)	
	4.5 U	UG/KG	Endrin	
	4.5 U	UG/KG	Endosulfan II (beta)	•
	4.5 U	UG/KG	4,4'-DDD (p,p'-DDD)	
	4.5 U	UG/KG	Endosulfan Sulfate	
	4.5 U	UG/KG	4,4'-DDT (ρ,ρ'-DDT)	
	23 U	UG/KG	Methoxychlor	
	4.5 U	UG/KG	Endrin Ketone	
	4.5 U	UG/KG	Endrin Aldehyde	
	2.3 U	UG/KG	alpha-Chlordane /2	•
`	2.3 U .	UG/KG	gamma-Chlordane /2	
	230 U	UG/KG	Toxaphene	
•	45 U	UG/KG	PCB-1016 (Aroclor 1016)	
	91 U	UG/KG	PCB-1221 (Aroclor 1221)	
	45 U	UG/KG	PCB-1232 (Aroclor 1232)	
	45 U	UG/KG	PCB-1242 (Aroclor 1242)	
	45 U	UG/KG	PCB-1248 (Aroclor 1248)	
	45 U	UG/KG	PCB-1254 (Aroclor 1254)	
	45 U	UG/KG	PCB-1260 (Aroclor 1260)	
	26	%	% Moisture	
			•	

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Produced by: Goddard. Denise Project: 03-0474 Sample 6796 FY 2003 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:50 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS06SB / MD No: 1XX6 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XX6 DATA REPORTED ON DRY WEIGHT BASIS

		<u>-</u>
RESULTS	UNITS	ANALYTE
2.3 U	UG/KG	alpha-BHC
2.3 U	UG/KG	beta-BHC
2.3 U	UG/KG	delta-BHC
2.3 U	UG/KG	gamma-BHC (Lindane)
2.3 U	UG/KG	Heptachlor
2.3 U	UG/KG	Aldrin
2.3 U	UG/KG	Heptachlor Epoxide
2.3 U	UG/KG	Endosulfan I (alpha)
4.5 U	UG/KG	Dieldrin
4.5 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.5 U 1	UG/KG	Endrin
4.5 U	UG/KG	Endosulfan II (beta)
4.5 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.5 U	UG/KG	Endosulfan Sulfate
4.5 U	UG/KG	4,4'-DDT (p,p'-DDT)
23 U	UG/KG	Methoxychlor
4.5 U	UG/KG	Endrin Ketone
4.5 U	UG/KG	Endrin Aldehyde
2.3 U	UG/KG	alpha-Chlordane /2
2.3 U	UG/KG	gamma-Chlordane /2
230 U	UG/KG	Toxaphene
45 U	UG/KG	PCB-1016 (Aroclor 1016)
91 U	UG/KG	PCB-1221 (Aroclor 1221)
45 U	UG/KG	PCB-1232 (Aroclor 1232)
45 U	UG/KG	PCB-1242 (Aroclor 1242)
45 U	UG/KG	PCB-1248 (Aroclor 1248)
45 U	UG/KG	PCB-1254 (Aroclor 1254)
45 U	UG/KG	PCB-1260 (Aroclor 1260)
26	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see:chlordane constituents | /2-constituents or metabolites of technical chlordane

Produced by: Goddard, Denise Project: 03-0474 FY 2003 Sample Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:35 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS24SS / MD No: 1XX7 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX7 DATA REPORTED ON DRY WEIGHT BASIS

	·	
RESULTS	UNITS	ANALYTE
3.1 U	UG/KG	alpha-BHC
5.3 U ·	UG/KG	beta-BHC
3.1 U	UG/KG	delta-BHC
3.1 U	UG/KG	gamma-BHC (Lindane)
3.1 U	UG/KG	Heptachlor
3.1 U	UG/KG	Aldrin
3.1 U	UG/KG	Heptachlor Epoxide
3.1 U	UG/KG	Endosulfan I (alpha)
6.0 U	UG/KG	Dieldrin
6.0 U	UG/KG.	4,4'-DDE (p,p'-DDE)
6.0 U	UG/KG	Endrin
6.0 U	UG/KG	Endosulfan II (beta)
6.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
6.0 U	UG/KG	Endosulfan Sulfate
6.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
31 U	UG/KG	Methoxychlor
6.0 U	UG/KG	Endrin Ketone
6.0 U	UG/KG	Endrin Aldehyde
3.1 U	UG/KG	alpha-Chiordane /2
3.1 U	UG/KG	gamma-Chlordane /2
310 U	UG/KG	Toxaphene
60 U	UG/KG	PCB-1016 (Aroclor 1016)
120 U	UG/KG	PCB-1221 (Aroclor 1221)
60 U	UG/KG	PCB-1232 (Aroclor 1232)
60 U	UG/KG	PCB-1242 (Aroclor 1242)
60 U	UG/KG	PCB-1248 (Aroclor 1248)
60 U	UG/KG	PCB-1254 (Aroclor 1254)
60 U	UG/KG	PCB-1260 (Aroclor 1260)
45 .	%	% Moisture
		į

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolities of technical chlordane

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 07:40 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS12SS / MD No: 1XX9 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX9 DATA REPORTED ON DRY WEIGHT BASIS

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RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
2.0 U	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.0 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
4.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
20 U	UG/KG	Methoxychlor
4.0 U	UG/KG	Endrin Ketone
4.0 U	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U .	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
40 U	UG/KG	PCB-1016 (Aroclor 1016)
81 U	UG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
40 U	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
17	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 07:50

DATA REPORTED ON DRY WEIGHT BASIS

Endina:

Media: SUBSURFACE SOIL (> 12")

6799 FY 2003

Case No: 31635 MD No: 1XY0

Flowood, MS

Inorg Contractor: SENTIN

D No: 1XY0

Org Contractor: LIBRTY

Facility: Gulf States Creosoting

Sample

**Pesticides Scan** 

Id/Station: GS12SB /

**RESULTS UNITS** 

Program: SF

2.3 U UG/KG alpha-BHC beta-BHC 2.3 U UG/KG delta-BHC 2.3 U UG/KG 2.3 U UG/KG gamma-BHC (Lindane)

**ANALYTE** 

2.3 U UG/KG Heptachlor 2.3°U UG/KG Aldrin

2.3 U UG/KG Heptachlor Epoxide 2.3 U UG/KG Endosulfan I (alpha)

4.5 U UG/KG Dieldrin

4.5 U 4,4'-DDE (p,p'-DDE) UG/KG

4.5 U UG/KG Endrin 4.5 U UG/KG Endosulfan II (beta) 4.5 U UG/KG 4.4'-DDD (p.p'-DDD) 4.5 U Endosulfan Sulfate UG/KG 4.5 U 4,4'-DDT (p,p'-DDT) UG/KG 23 U UG/KG Methoxychlor 4.5 U UG/KG Endrin Ketone 4.5 U UG/KG Endrin Aldehyde 2.3 U UG/KG alpha-Chlordane /2

45 U UG/KG PCB-1016 (Aroclor 1016) 91 U UG/KG PCB-1221 (Aroclor 1221)

Toxaphene

gamma-Chlordane /2

PCB-1260 (Arocior 1260)

45 U UG/KG PCB-1232 (Aroclor 1232) 45 U UG/KG PCB-1242 (Aroclor 1242) 45 U UG/KG PCB-1248 (Aroclor 1248) PCB-1254 (Aroclor 1254) 45 U UG/KG

% Moisture

45 U UG/KG 26 %

UG/KG

UG/KG

2.3 U

230 U

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Produced by: Goddard, Denise Project: 03-0474 Sample FY 2003 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 07:40 Program: SF Case No: 31635 Ending: Id/Station: GS03SS / Inorg Contractor: SENTIN MD No: 1XY1 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XY1 DATA REPORTED ON DRY WEIGHT BASIS R

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.3 U	UG/KG	Dieldrin
4.3 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.3 U	UG/KG	Endrin
4.3 U	UG/KG	Endosulfan II (beta)
4.3 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.3 U	UG/KG	Endosulfan Sulfate
4.3 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.3 U	UG/KG	Endrin Ketone
. 4.3 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
220 U	UG/KG	Toxaphene
43 U	UG/KG	PCB-1016 (Aroclor 1016)
88 U	UG/KG	PCB-1221 (Aroctor 1221)
43 U	UG/KG	PCB-1232 (Aroclor 1232)
43 U	UG/KG	PCB-1242 (Aroclor 1242)
43 U	UG/KG	PCB-1248 (Aroclor 1248)
43 U	UG/KG	PCB-1254 (Aroclor 1254)
43 U	UG/KG	PCB-1260 (Arocior 1260)
24	%	% Moisture
•		•

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Sample FY 2003 Project: 03-0474

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS03SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XY2 D No: 1XY2

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 07:55

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.4 U	UG/KG	alpha-BHC
2.4 U	UG/KG	beta-BHC
2.4 U	UG/KG	delta-BHC
2.4 U	UG/KG	gamma-BHC (Lindane)
2.4 U	UG/KG	Heptachlor
2.4 U	UG/KG	Aldrin
2.4 U	UG/KG	Heptachlor Epoxide
2.4 U	UG/KG	Endosulfan I (alpha)
4.6 U	UG/KG	Dieldrin
4.6 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.6 U	UG/KG	Endrin
4.6 U	UG/KG	Endosulfan II (beta)
4.6 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.6 U	UG/KG	Endosulfan Sulfate
4.6 U	UG/KG	4,4'-DDT (p,p'-DDT)
24 U	UG/KG	Methoxychlor
4.6 U	UG/KG	Endrin Ketone
4.6 U	UG/KG	Endrin Aldehyde
2.4 U	UG/KG	alpha-Chlordane /2
2.4 U	UG/KG	gamma-Chlordane /2
240 U ·	UG/KG	Toxaphene
46 U	UG/KG	PCB-1016 (Aroclor 1016)
93 U	UG/KG	PCB-1221 (Aroclor 1221)
46 U	UG/KG	PCB-1232 (Aroclor 1232)
46 U	UG/KG	PCB-1242 (Aroclor 1242)
46 U	UG/KG	PCB-1248 (Aroclor 1248)
46 U	UG/KG	PCB-1254 (Aroclor 1254)
46 U	UG/KG	PCB-1260 (Aroclor 1260)
28	%	% Moisture
•		·

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

FY 2003 Project: 03-0474 Sample 6802

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS08SS /

Media: SURFACE SOIL (0" - 12")

Case No: 31635

MD No: 1XY3

D No: 1XY3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:10

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
1.8 U	UG/KG	alpha-BHC
1.8 U	UG/KG	beta-BHC
1.8 U	UG/KG	delta-BHC
1.8 U	UG/KG	gamma-BHC (Lindane)
1.8 U	UG/KG	Heptachlor
1.8 U	UG/KG	Aldrin
1.8 U	UG/KG	Heptachlor Epoxide
1.8 U	UG/KG	Endosulfan I (alpha)
3.4 U	UG/KG	Dieldrin
13 U	UG/KG	4,4'-DDE (p,p'-DDE)
10 U	UG/KG	Endrin
3.4 U •	UG/KG	Endosulfan II (beta)
3.4 U	UG/KG	4,4'-DDD (p,p'-DDD)
3.4 U	UG/KG	Endosulfan Sulfate
41 J	UG/KG	4,4'-DDT (p,p'-DDT)
74 NJ	UG/KG	Methoxychlor
24 U	UG/KG	Endrin Ketone
17 J	UG/KG	Endrin Aldehyde
1.8 U	UG/KG	alpha-Chlordane /2
1.8 U	UG/KG	gamma-Chlordane /2
180 U	UG/KG	Toxaphene
34 U	UG/KG	PCB-1016 (Aroclor 1016)
70 U	UG/KG	PCB-1221 (Aroclor 1221)
34 U	UG/KG	PCB-1232 (Aroclor 1232)
34 U	UG/KG	PCB-1242 (Aroclor 1242)
34 U	UG/KG	PCB-1248 (Aroclor 1248)
34 U	UG/KG	PCB-1254 (Aroclor 1254)
34 U	UG/KG	PCB-1260 (Aroclor 1260)
4	%	% Moisture
		İ

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

Flowood, MS

Production Date: 06/20/2003 14:11

Project: 03-0474 Sample 6804 FY 2003

Produced by: Goddard, Denise

**Pesticides Scan** 

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:20

Program: SF

Ending:

ld/Station: GS08SB /

Case No: 31635 MD No: 1XY4

Media: SUBSURFACE SOIL (> 12")

Inorg Contractor: SENTIN

Org Contractor: LIBRTY D No: 1XY4

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U ·	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
· 2.0 U	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
4.0 U	UG/KG	Dieldrin
4.0 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.0 U	UG/KG	Endrin
4.0 U	UG/KG	Endosulfan II (beta)
4.0 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.0 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG-	4,4'-DDT (p,p'-DDT)
20 U	UG/KG	Methoxychlor
4.0 U	UG/KG	Endrin Ketone
4.0 U	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
40 U	UG/KG	PCB-1016 (Aroclor 1016)
81 U	UG/KG	PCB-1221 (Aroclor 1221)
40 U	UG/KG	PCB-1232 (Aroclor 1232)
40 U	UG/KG	PCB-1242 (Aroclor 1242)
40 U	UG/KG	PCB-1248 (Aroclor 1248)
40 U	UG/KG	PCB-1254 (Aroclor 1254)
40 U	UG/KG	PCB-1260 (Aroclor 1260)
17	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/20/2003 14:11

6805 . FY 2003 Project: 03-0474 Sample

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS07SS /

Media: SURFACE SOIL (0" - 12")

Case No: 31635

MD No: 1XY5 D No: 1XY5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:35

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	UG/KG	gamma-BHC (Lindane)
. 2.0 U	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
3.9 U	UG/KG	Dieldrin
6.4 U	UG/KG	4,4'-DDE (p,p'-DDE)
7.1 U	UG/KG	Endrin
3.9 U	UG/KG	Endosulfan II (beta)
3.9 U	UG/KG	4,4'-DDD (p,p'-DDD)
3.9 U	UG/KG	Endosulfan Sulfate
12 J	UG/KG	4,4'-DDT (p,p'-DDT)
47 U	UG/KG	Methoxychlor
20 U	UG/KG	Endrin Ketone
3.9 U	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
39 U -	UG/KG	PCB-1016 (Aroclor 1016)
79 U	UG/KG	PCB-1221 (Aroclor 1221)
39 U	UG/KG	PCB-1232 (Aroclor 1232)
39 U	UG/KG	PCB-1242 (Aroclor 1242)
39 U	UG/KG	PCB-1248 (Aroclor 1248)
39 U	UG/KG	PCB-1254 (Aroclor 1254)
39 U	UG/KG	PCB-1260 (Arocior 1260)
15	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

PESTICIDES/PCB SAMPLE ANALYSIS	EPA - REGION IV SESD
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Project: 03-0474

D. ATHENS, GA

Production Date: 06/20/2003 14:11

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:45

Endina:

Case No: 31635 Inorg Contractor: SENTIN MD No: 1XY6

Id/Station: GS07SB /

Flowood, MS

Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XY6 DATA REPORTED ON DRY WEIGHT BASIS

**RESULTS UNITS ANALYTE** 2.3 U alpha-BHC UG/KG beta-BHC 2.3 U UG/KG 2.3 U UG/KG delta-BHC 2.3 U UG/KG gamma-BHC (Lindane) 2.3 U Heptachlor UG/KG 2.3 U UG/KG Aldrin 2.3 U Heptachlor Epoxide UG/KG 2.3 U Endosulfan I (alpha) UG/KG 4.5 U UG/KG Dieldrin 4.4'-DDE (p,p'-DDE) 4.5 U UG/KG 4.5 U Endrin UG/KG 4.5 U UG/KG Endosulfan II (beta) 4.5 U UG/KG 4.4'-DDD (p.p'-DDD) 4.5 U UG/KG Endosulfan Sulfate 4,4'-DDT (p.p'-DDT) 4.5 U UG/KG 23 U UG/KG Methoxychlor 4.5 U UG/KG **Endrin Ketone** 4.5 U UG/KG Endrin Aldehyde 2.3 U UG/KG alpha-Chlordane /2 gamma-Chlordane /2 2.3 U UG/KG 230 U UG/KG Toxaphene 45 U PCB-1016 (Aroclor 1016) UG/KG PCB-1221 (Aroclor 1221) 91 U UG/KG PCB-1232 (Aroclor 1232) 45 U UG/KG PCB-1242 (Aroclor 1242) 45 U UG/KG 45 U UG/KG PCB-1248 (Aroclor 1248) 45 U PCB-1254 (Aroclor 1254) UG/KG 45 U UG/KG PCB-1260 (Aroclor 1260) 26 % Moisture

6806 FY 2003

Facility: Gulf States Creosoting

Sample

**Pesticides Scan** 

Program: SF

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Produced by: Goddard, Denise Sample FY 2003 Project: 03-0474 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 08:30 Program: SF Case No: 31635 Ending: Id/Station: GS04SS / Inorg Contractor: SENTIN MD No: 1XY7 **Org Contractor: LIBRTY** Media: SURFACE SOIL (0" - 12") D No: 1XY7 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 ∪•	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KĠ	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.3 U	UG/KG	Dieldrin
4.3 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.3 U	UG/KG	Endrin
4.3 U	UG/KG	Endosulfan II (beta)
4.3 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.3 U	UG/KG	Endosulfan Sulfate
4.3 U `	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.3 U	UG/KG	Endrin Ketone
4.3 U	UG/KG	Endrin Aldehyde
- 2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
. 220 U	UG/KG	Toxaphene
43 U	UG/KG	PCB-1016 (Aroclor 1016)
88 U	UG/KG	PCB-1221 (Aroclor 1221)
43 U	UG/KG	PCB-1232 (Aroclor 1232)
43 U	UG/KG	PCB-1242 (Aroclor 1242)
43 U	UG/KG	PCB-1248 (Aroclor 1248)
43 U	UG/KG	PCB-1254 (Aroclor 1254)
43 U	UG/KG	PCB-1260 (Aroclor 1260)
24	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value. NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474 Sample 6808 FY 2003

Pesticides Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS04SB /

Case No: 31635

Media: SUBSURFACE SOIL (> 12")

MD No: 1XY8 D No: 1XY8

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2.2 U	UG/KG	alpha-BHC
2.2 U	UG/KG	beta-BHC
2.2 U	UG/KG	delta-BHC
2.2 U	UG/KG	gamma-BHC (Lindane)
2.2 U	UG/KG	Heptachlor
2.2 U	UG/KG	Aldrin
2.2 U	UG/KG	Heptachlor Epoxide
2.2 U	UG/KG	Endosulfan I (alpha)
4.3 U	UG/KG	Dieldrin
4.3 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.3 U	UG/KG	Endrin
4.3 U	UG/KG	Endosulfan II (beta)
4.3 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.3 U	UG/KG	Endosulfan Sulfate
4.3 U	UG/KG	4,4'-DDT (p,p'-DDT)
22 U	UG/KG	Methoxychlor
4.3 U	UG/KG	Endrin Ketone
4.3 U	UG/KG	Endrin Aldehyde
2.2 U	UG/KG	alpha-Chlordane /2
2.2 U	UG/KG	gamma-Chlordane /2
.220 U	UG/KG	Toxaphene
43 U •	UG/KG	PCB-1016 (Aroclor 1016)
87 U	UG/KG	PCB-1221 (Aroclor 1221)
43 U	UG/KG	PCB-1232 (Aroclor 1232)
43 U	UG/KG	PCB-1242 (Aroclor 1242)
43 U	UG/KG	PCB-1248 (Aroclor 1248)
43 U	UG/KG	PCB-1254 (Aroclor 1254)
43 U	UG/KG	PCB-1260 (Aroclor 1260)
23	%	% Moisture
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U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Produced by: Goddard, Denise Project: 03-0474 Sample 6809 FY 2003 Requestor: **Pesticides Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 09:15 Program: SF Case No: 31635 Ending: Id/Station: GS05SS / Inorg Contractor: SENTIN MD No: 1XY9 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XY9 DATA REPORTED ON DRY WEIGHT BASIS

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RESULTS	UNITS	ANALYTE
2.0 U	UG/KG	alpha-BHC
2.0 U	UG/KG	beta-BHC
2.0 U	UG/KG	delta-BHC
2.0 U	.UG/KG	gamma-BHC (Lindane)
2.0 U	UG/KG	Heptachlor
2.0 U	UG/KG	Aldrin
. 2.0 U	UG/KG	Heptachlor Epoxide
2.0 U	UG/KG	Endosulfan I (alpha)
3.9 U	UG/KG	Dieldrin
9.9 U	UG/KG	4,4'-DDE (p,p'-DDE)
12 U	UG/KG	Endrin
3.9 U	UG/KG	Endosulfan II (beta)
3.9 U .	UG/KG	4,4'-DDD (p,p'-DDD)
3.9 U	UG/KG	Endosulfan Sulfate
4.0 U	UG/KG	4,4'-DDT (p,p'-DDT)
48 U	UG/KG	Methoxychlor
17 U	UG/KG	Endrin Ketone
18 J	UG/KG	Endrin Aldehyde
2.0 U	UG/KG	alpha-Chlordane /2
2.0 U	UG/KG	gamma-Chlordane /2
200 U	UG/KG	Toxaphene
39 U	UG/KG	PCB-1016 (Aroclor 1016)
79 U	UG/KG	PCB-1221 (Aroclor 1221)
39 U	UG/KG	PCB-1232 (Aroclor 1232)
39 U	UG/KG	PCB-1242 (Aroclor 1242)
39 U	UG/KG	PCB-1248 (Aroclor 1248)
39 U	UG/KG	PCB-1254 (Aroclor 1254)
39 U	UG/KG	PCB-1260 (Aroclor 1260)
15	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474 Sample 6810 FY 2003

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS05SB /

MD No: 1XZ0

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

D No: 1XZ0

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:25

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

ESULTS	UNITS	ANALYTE
2.1 U	UG/KG	alpha-BHC
2.1 U	UG/KG	beta-BHC
2.1 U	UG/KG	delta-BHC
2.1 U -	UG/KG	gamma-BHC (Lindane)
2.1 U	UG/KG	Heptachlor
2.1 U	UG/KG	Aldrin
2.1 U	UG/KG	Heptachlor Epoxide
2.1 U	UG/KG	Endosulfan I (alpha)
4.1 U	UG/KG	Dieldrin
4.1 U	UG/KG	4,4'-DDE (p,p'-DDE)
4.1 U	UG/KG	Endrin
4.1 U	UG/KG	Endosulfan II (beta)
4.1 U	UG/KG	4,4'-DDD (p,p'-DDD)
4.1 U	UG/KG	Endosulfan Sulfate
4.1 U	UG/KG	4,4'-DDT (p,p'-DDT)
21 U	UG/KG	Methoxychlor
4.1 U	UG/KG	Endrin Ketone
4.1 U	UG/KG	Endrin Aldehyde
2.1 U	UG/KG	alpha-Chlordane /2
2.1 U	UG/KG	gamma-Chlordane /2
210 U	UG/KG	Toxaphene
41 U	UG/KG	PCB-1016 (Aroclor 1016)
83 U	UG/KG	PCB-1221 (Aroclor 1221)
41 U	UG/KG	PCB-1232 (Aroclor 1232)
41 U	UG/KG	PCB-1242 (Aroclor 1242)
41 U	UG/KG	PCB-1248 (Aroclor 1248)
41 U	UG/KG	PCB-1254 (Arocior 1254)
41 U	UG/KG	PCB-1260 (Aroclor 1260)
19 ,	%	% Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Sample

6811

Project: 03-0474 FY 2003

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS03GW /

MD No: 1XZ1

Media: GROUNDWATER

Case No: 31635

D No: 1XZ1

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:55

Ending:

RESULTS	UNITS	ANALYTE
0.010 U	UG/L	alpha-BHC
0.010 U	UG/L	beta-BHC
0.010 U	UG/L	delta-BHC
0.010 U	UG/L	gamma-BHC (Lindane)
0.010 U	UG/L	Heptachlor
0.010 U	UG/L	Aldrin
0.010 U	UG/L	Heptachlor Epoxide
0.010 U	UG/L	Endosulfan I (alpha)
0.020 U	UG/L	Dieldrin
0.020 UJ -	UG/L	4,4'-DDE (p,p'-DDE)
0.020 UJ	UG/L	Endrin
0.020 U	UG/L	Endosulfan II (beta)
0.020 UJ	UG/L	4,4'-DDD (p,p'-DDD)
0.020 U	UG/L	Endosulfan Sulfate
0.020 UJ	UG/L	4,4'-DDT (p,p'-DDT)
0.10 U	UG/L	Methoxychlor
0.020 UJ	UG/L	Endrin Ketone
0.020 UJ	UG/L	Endrin Aldehyde
0.010 U	UG/L	alpha-Chlordane /2
0.010 U	UG/L	gamma-Chlordane /2
1.0 U	UG/L	Toxaphene
0.20 U	UG/L	PCB-1016 (Aroclor 1016)
0.40 U	UG/L	PCB-1221 (Aroclor 1221)
0.20 U	UG/L	PCB-1232 (Aroclor 1232)
0.20 U	UG/L	PCB-1242 (Aroclor 1242)
0.20 U	UG/L	PCB-1248 (Aroclor 1248)
0.20 U	UG/L	PCB-1254 (Aroclor 1254)
0.20 U	UG/L	PCB-1260 (Aroclor 1260)

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

FY 2003 Project: 03-0474 Sample 6812

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flawood, MS

Program: SF

Case No: 31635

Id/Station: GS04GW / MD No: 1XZ3

Media: GROUNDWATER

D No: 1XZ3

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 13:20

Ending:

RESULTS	UNITS	ANALYTE
0.010 U	UG/L	alpha-BHC
0.010 U	UG/L	beta-BHC
0.010 U	UG/L	delta-BHC
0.010 U	'UG/L	gamma-BHC (Lindane)
0.010 U	UG/L	Heptachlor
0.010 U	UG/L	Aldrin
0.010 U	UG/L	Heptachlor Epoxide
0.010 U	UG/L	Endosulfan I (alpha)
0.020 U	UG/L	Dieldrin
0.020 UJ	UG/L	4,4'-DDE (p,p'-DDE)
0.020 UJ	UG/L	Endrin
0.020 U	UG/L	Endosulfan II (beta)
0.020 UJ	UG/L	4,4'-DDD (p,p'-DDD)
0.020 U	UG/L	Endosulfan Sulfate
0.020 UJ	UG/L	4,4'-DDT (p,p'-DDT)
0.10 U	UG/L	Methoxychlor
0.020 UJ	UG/L	Endrin Ketone
0.020 UJ	UG/L	Endrin Aldehyde
0.010 U	UG/L	alpha-Chlordane /2
0.010 U	UG/L	gamma-Chlordane /2
1.0 U	UG/L	Toxaphene
0.20 U	UG/L	PCB-1016 (Aroclor 1016)
0.40 U	UG/L	PCB-1221 (Aroclor 1221)
0.20 U	UG/L	PCB-1232 (Aroclor 1232)
0.20 U	UG/L	PCB-1242 (Aroclor 1242)
0.20 U	UG/L	PCB-1248 (Aroclor 1248)
0.20 U	UG/L	PCB-1254 (Aroclor 1254)
0.20 U	UG/L	PCB-1260 (Aroclor 1260)
•		; i

U-Analyte not detected at or above reporting limit. | J-Idenţification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

6813 FY 2003 Project: 03-0474 Sample

**Pesticides Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS05GW / Media: GROUNDWATER Case No: 31635

MD No: 1XZ4 D No: 1XZ4

Inorg Contractor: SENTIN

Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 11:45

Ending:

UNITS	ANALYTE
UG/L	alpha-BHC
UG/L	beta-BHC
UG/L	delta-BHC
UG/L	gamma-BHC (Lindane)
UG/L	Heptachlor
UG/L	Aldrin
UG/L	Heptachlor Epoxide
UG/L	Endosulfan I (alpha)
UG/L	Dieldrin
UG/L	4,4'-DDE (p,p'-DDE)
UG/L	Endrin
UG/L	Endosulfan II (beta)
	4,4'-DDD (p,p'-DDD)
	Endosulfan Sulfate
	4,4'-DDT (p,p'-DDT)
	Methoxychlor
	Endrin Ketone
	Endrin Aldehyde
	alpha-Chlordane /2
	gamma-Chlordane /2
	Toxaphene
	PCB-1016 (Aroclor 1016)
	PCB-1221 (Aroclor 1221)
	PCB-1232 (Aroclor 1232)
	PCB-1242 (Aroclor 1242)
	PCB-1248 (Aroclor 1248)
	PCB-1254 (Aroclor 1254)
UG/L	PCB-1260 (Aroclor 1260)
	<u> </u>
	UG/L UG/L UG/L UG/L UG/L UG/L UG/L UG/L

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable. C-confirmed by GCMS | /1-when no value is reported, see chlordane constituents | /2-constituents or metabolites of technical chlordane

May 21, 2003

## INORGANIC DATA QUALIFIERS REPORT

Case Number: 31635 Project Number: 03-0474

Sample No	Element	Flaq	Reason
6754	Al	J	Serial dilution percent difference = 10.8%
	Sb	Ĵ	Matrix spike recovery = 45.5%
	Cu	Ū	Baseline instability in cal blanks
	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	Ni	Ü	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
	Se	R	Analyte reported as potential false positive
	Ag	R	Analyte reported as potential false positive
6755	Al	J	Serial dilution percent difference = 10.8%
	Sb	J	Matrix spike recovery = 45.5%
	Cu	U	Baseline instability in cal blanks
	Pb .	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	Ni	· U	Baseline instability in cal blanks
	K	Ĵ	Serial dilution percent difference = 42%
	Ag	R	Analyte reported as potential false positive
	_		
6756	Al	J	Serial dilution percent difference = 10.8%
	Sb	UJ	Matrix spike recovery = 45.5%
•		•	Baseline instability in cal blanks
	Cu	U	Baseline instability in cal blanks
	. Pb	J	Matrix spike recovery = 184.2%
	Mg	.J	Serial dilution percent difference = 14.1%
	Ni	U	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
6757	· Al	J	Serial dilution percent difference = 10.8%
6757	Sb	ບັນ	Matrix spike recovery = 45.5%
	SD		Baseline instability in cal blanks
	C) v	U	Baseline instability in cal blanks
	Cu		
•	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	, Ni	Ŭ	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
	Se	J	% RSD > 20% for ICP multiple exposures
6758	Al .	J	Serial dilution percent difference = 10.8%
0,50	Sb	UJ .	Matrix spike recovery = 45.5%
	-~	,==	Baseline instability in cal blanks
	Cu	ΰ	Baseline instability in cal blanks
	Pb	J	Matrix spike recovery = 184.2%
		J	
	Mg	υ. υ	
	Hg		Positive reported < lowest std on cal curve
	Ni	U	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
	Ag	R	Analyte reported as potential false positive

# **UNITED STATES ENVIRONMENTAL PROTECTION AGENCY**

#### Region 4

Science and Ecosystem Support Division 980 College Station Road Athens, Georgia 30605-2720

#### **MEMORANDUM**

Date: 06/05/2003

Subject: Results of METALS Sample Analysis

03-0474

**Gulf States Creosoting** 

Flowood, MS

From: Goddard, Denise

\*To: Striggow, Brian

Thru: QA Office

Attached are the results of analysis of samples collected as part of the subject project. If you have any questions, please contact me.

May 21, 2003

## INORGANIC DATA QUALIFIERS REPORT (continued)

Case Number: 31635

Project Number: 03-0474

Sample No.	Element	Flag	Reason
6759	Al	J.	Serial dilution percent difference = 10.8%
	Sb	UĴ	Matrix spike recovery ≈ 45.5%
			Baseline instability in cal blanks
	As	<sup>'</sup> R	Analyte reported as potential false positive
	Cu	U	Baseline instability in cal blanks
,	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
•	Ni	Ŭ.	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
	Ag	R	Analyte reported as potential false positive
6760	Al	. <b>J</b>	Serial dilution percent difference = 10.8%
	Sb	UJ	Matrix spike recovery = 45.5%
•		V.	Baseline instability in cal blanks
	Cd	·R	Analyte reported as potential false positive
	Cu	U	Baseline instability in cal blanks
,	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	Нg	U	Positive reported < lowest std on cal curve
	Ni	U	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
	Se	J	% RSD > 20% for ICP multiple exposures
•	Ag	R	Analyte reported as potential false positive
6761	Al .	• Ј	Serial dilution percent difference = 10.8%
	Sb	UJ	Matrix spike recovery = 45.5%
			Baseline instability in cal blanks
	Cu	U	Baseline instability in cal blanks
•	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
•	Ni	U .	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
	Ag	R	Analyte reported as potential false positive
6762	Al	J	Serial dilution percent difference = 10.8%
	Sb	J	Matrix spike recovery = 45.5%
	Cu		Baseline instability in cal blanks
•	Pb	J,	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
•	Нg	Ū	Positive reported < lowest std on cal curve
•	Ni	ប	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
·.	Se	J	% RSD > 20% for ICP multiple exposures
	Ag	R	Analyte reported as potential false positive

May 21, 2003

## INORGANIC DATA QUALIFIERS REPORT (continued)

Case Number: 31635

Project Number: 03-0474

Sample	No.	Element	Flaq	Reason
6763		Al	J	Serial dilution percent difference = 10.8%
·		Sb	J	Matrix spike recovery = 45.5%
•	•	Cu	ับ	Baseline instability in cal blanks
		Pb	J	Matrix spike recovery = 184.2%
		Mg	J	Serial dilution percent difference = 14.1%
		Ni	ָּט ·	Baseline instability in cal blanks
		K	J	Serial dilution percent difference = 42%
	•	Ag	R	Analyte reported as potential false positive
6764		Al	J	Serial dilution percent difference = 10.8%
0,01		Sb	J	Matrix spike recovery = 45.5%
		Cu	บ	Baseline instability in cal blanks
		Pb	J	Matrix spike recovery = 184.2%
		Mg	ับ	Serial dilution percent difference = 14.1%
•				
		Hg	U	Positive reported < lowest std on cal curve
		Ni	Ŭ	Baseline instability in cal blanks
		K	J	Serial dilution percent difference = 42%
		Se	R	Analyte reported as potential false positive
6765		Al	J	Serial dilution percent difference = 10.8%
		Sb	· UJ	Matrix spike recovery = 45.5%
				Baseline instability in cal blanks
		As	R	Analyte reported as potential false positive
		Cu	U	Baseline instability in cal blanks
		Pb	J	Matrix spike recovery = 184.2%
		Mg	· J	Serial dilution percent difference = 14.1%
		Ni	Ų	Baseline instability in cal blanks
·		K	Ĵ	Serial dilution percent difference = 42%
		Ag	R	Analyte reported as potential false positive
6766		Al	J	Serial dilution percent difference = 10.8%
		Sb	UJ	Matrix spike recovery = 45.5%
. '				Baseline instability in cal blanks
		As	J	% RSD > 20% for ICP multiple exposures
		Cu	Ū	Baseline instability in cal blanks
·				Matrix spike recovery = 184.2%
		Mg	J	Serial dilution percent difference = 14.1%
		Ni	Ū	Baseline instability in cal blanks
		K	, J	Serial dilution percent difference = 42%
		Ag	R	Analyte reported as potential false positive
6767		Al	J	Serial dilution percent difference = 10.8%
		Sb	J	Matrix spike recovery = 45.5%
		Pb ·	J	Matrix spike recovery = 184.2%
•		Mg	J	Serial dilution percent difference = 14.1%
		Ni	Ü	Baseline instability in cal blanks
		K	J	Serial dilution percent difference = 42%
•		Se	J	% RSD > 20% for ICP multiple exposures
				· · · · · · · · · · · · · · · · · · ·

May 21, 2003 INORGANIC DATA QUALIFIERS REPORT (continued)

Case Number: 31635 Project Number: 03-0474

Sample No.	Element	Flaq_	Reason
6768	Al	J	Serial dilution percent difference = 10.8%
	Sb	UJ	Matrix spike recovery = 45.5%
			Baseline instability in cal blanks
	Cu	U	Baseline instability in cal blanks
	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	Ni	U	Baseline instability in cal blanks
•	K	J	Serial dilution percent difference = 42%
6769	Al	` J	Serial dilution percent difference = 10.8%
• ,	Sb	J	Matrix spike recovery = 45.5%
	Cu	U	Baseline instability in cal blanks
	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	Ni	U	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
•	Se	R	Analyte reported as potential false positive
3770	Al	IJ	PE sample recovery > warning limit
		•	PE sample recovery > action limit
			Positive in blind blank
	Cr	R	Analyte reported as potential false positive
	Cu	. <b>U</b>	Baseline instability in blind blank
6771	Al	J	Serial dilution percent difference = 10.8%
0772	Sb	J	Matrix spike recovery = 45.5%
•	. Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	Ni	ับ	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
6772	Al	J	Serial dilution percent difference = 10.8%
5772	Sb	J	Matrix spike recovery = 45.5%
	Pb	J	Matrix spike recovery = 184.2%
	Mg	J	Serial dilution percent difference = 14.1%
	Ni	Ū	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42%
	Ag	R	Analyte reported as potential false positive
6773	Al	· J	Serial dilution percent difference = 13.4%
	Sb	R	Matrix spike recovery = 50.8%
			Analyte reported as potential false positive
	€đ	·R	Analyte reported as potential false positive
	Cu	Ĵ	Serial dilution percent difference = 45.6%
	Mg	Ĵ	Serial dilution percent difference = 14.9%
	Hg	ŭ	Positive reported < lowest std on cal curve
	Ni	ັນ .	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
	Ag.	R	Analyte reported as potential false positive
·			•

May 21, 2003 INORGANIC DATA QUALIFIERS REPORT (continued)

Project Number: 03-0474

Sample No.	Element	Flag	Reason
6774	Al'	. J	Serial dilution percent difference = 13.4%
	Sb	R	Matrix spike recovery = 50.8%
		•	Analyte reported as potential false positive
•	Cu	ŪJ	Serial dilution percent difference = 45.6%
			Baseline instability in prep blank
	Mg	J	Serial dilution percent difference = 14.9%
	Ni	. Т	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
6775	Al	J	Serial dilution percent difference = 13.4%
<b>V</b> / · · ·	Sb	Ĵ	Matrix spike recovery = 50.8%
	Cu	บัง	Serial dilution percent difference = 45.6%
		-	Baseline instability in prep blank
	Mg	J	Serial dilution percent difference = 14.9%
	Ni	บ	Baseline instability in cal blanks
	K .	J	Serial dilution percent difference = 42.7%
	Ag	R	Analyte reported as potential false positive
6776	Al	J	Coming dilution annual differences 12 At
6776	Sb		Serial dilution percent difference = 13.4%
	20	R	Matrix spike recovery = 50.8%
	λ σ	70	Analyte reported as potential false positive
•	As	R	Analyte reported as potential false positive
	Cu	ປັ	Serial dilution percent difference = 45.6%
	Ma	7	Baseline instability in prep blank
	Mg	J	Serial dilution percent difference = 14.9%
	Ni	Ŭ	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
	Ag	R	Analyte reported as potential false positive
6777	Al	J	Serial dilution percent difference = 13.4%
	Sb	J	Matrix spike recovery = 50.8%
			Analyte reported as potential false positive
	Cr	U	Baseline instability in cal blanks
	Cu	ŬĴ	Serial dilution percent difference = 45.6%
			Baseline instability in prep blank
	Pb	J	% RSD > 20% for ICP multiple exposures
	Mg		Serial dilution percent difference = 14.9%
	Ni	U	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
	V	R	Analyte reported as potential false positive
6778	Al	J	Serial dilution percent difference = 13.4%
	Sb	J.	Matrix spike recovery = 50.8%
i	Cd	R	Analyte reported as potential false positive
	Cu	· J	Serial dilution percent difference = 45.6%
	Mg	·J	Serial dilution percent difference = 14.9%
•	Нg	Ū	Positive reported < lowest std on cal curve
•	Ni	Ŭ	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
	Se	. R	Analyte reported as potential false positive
•	Ag	R	Analyte reported as potential false positive
			1 F The Foot

May 21, 2003 INORGANIC DATA QUALIFIERS REPORT (continued)

Project Number: 03-0474

_		_		
<u>Sample</u>	No.	Element	Flag	Reason
6779		Al	J	Serial dilution percent difference = 13.4%
		Sb	J	Matrix spike recovery = 50.8%
		Cu	びび	Serial dilution percent difference = 45.6%
				Baseline instability in prep blank
		Mg	J	Serial dilution percent difference = 14.9%
•		Ni	Ü	Baseline instability in cal blanks
	•	K	Ĵ	Serial dilution percent difference = 42.7%
		Se	R	Analyte reported as potential false positive
		JC	IX.	Midiyee reported as potential rathe positive
6780		Al	J	Serial dilution percent difference = 13.4%
6780		Sb	J	Matrix spike recovery = 50.8%
			J	
		Cu		•
	•	Mg	J 	Serial dilution percent difference = 14.9%
		Ni	Ŭ	Baseline instability in cal blanks
		K	J	Serial dilution percent difference = 42.7%
		. Se	J	% RSD > 20% for ICP multiple exposures
•		- •	· _	- 1.2. 113
6781		Al	J	Serial dilution percent difference = 13.4%
•	•	Sb	R	Matrix spike recovery = 50.8%
				Analyte reported as potential false positive
•		As	J	% RSD > 20% for ICP multiple exposures
		Cu	${f U}{f J}$	Serial dilution percent difference = 45.6%
		•		Baseline instability in prep blank
		Mg	· J	Serial dilution percent difference = 14.9%
		Νī	U	Baseline instability in cal blanks
		ĸ	J	Serial dilution percent difference = 42.7%
		Se	R	Analyte reported as potential false positive
		. Ag	R	Analyte reported as potential false positive
		_		
6782		Al	J	Serial dilution percent difference = 13.4%
	• •	Sb	J	Matrix spike recovery = 50.8%
		As	J	% RSD > 20% for ICP multiple exposures
		Cu	UJ	Serial dilution percent difference = 45.6%
•				Baseline instability in prep blank
		Mg	J	Serial dilution percent difference = 14.9%
		Ni	Ū	Baseline instability in cal blanks
		K-:: ·	<del>"</del>	Serial dilution percent difference = 42.7%
		Ag	R	Analyte reported as potential false positive
		**9	•	initiate reported up possibility rather positive
6783		Al	J	Serial dilution percent difference = 13.4%
• •		Sb	J	Matrix spike recovery = 50.8%
		cđ	R	Analyte reported as potential false positive
		Cu	ับ บั	Serial dilution percent difference = 45.6%
		Cu		Baseline instability in prep blank
		Ma	т.	Serial dilution percent difference = 14.9%
		Mg	J	
		Hg	U	Positive reported < lowest std on cal curve
		Ni	ū	Baseline instability in cal blanks
		K	J	Serial dilution percent difference = 42.7%
		Se	R	Analyte reported as potential false positive

May 21, 2003 INORGANIC DATA QUALIFIERS REPORT (continued)

Project Number: 03-0474

Sample No.	Element	Flag	Reason
6784	Al	J	Serial dilution percent difference = 13.4%
	Sb	J	Matrix spike recovery = 50.8%
	Cu	UJ	Serial dilution percent difference = 45.6%
		•	Baseline instability in prep blank
	Mg	J	Serial dilution percent difference = 14.9%
	Ni	Ū	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
6785	Al	J	Serial dilution percent difference = 13.4%
	Sb	R	Matrix spike recovery = 50.8%
			Analyte reported as potential false positive
	Cu	J	Serial dilution percent difference = 45.6%
•	Mg	' J	Serial dilution percent difference = 14.9%
	нg	บ	Positive reported < lowest std on cal curve
	Νĭ	บ	Baseline instability in cal blanks
	К	J	Serial dilution percent difference = 42.7%
	. Se	R ·	Analyte reported as potential false positive
6786	Al	J <sup>.</sup>	Serial dilution percent difference = 13.4%
	Sb	R	Matrix spike recovery = 50.8%
			Analyte reported as potential false positive
	Cu	· J	Serial dilution percent difference = 45.6%
	Mg	J	Serial dilution percent difference = 14.9%
	Нg	บ	Positives reported < lowest std on cal curve
	ĸ	J	Serial dilution percent difference = 42.7%
3787	Al	J	PE sample recovery > warning limit
			PE sample recovery > action limit
	Co	R	Analyte reported as potential false positive
	Cu	U	Baseline instability in blind blank
	V	R	Analyte reported as potential false positive
6788	Al	J	Serial dilution percent difference = 13.4%
*)	Sb	J	Matrix spike recovery = 50.8%
	Cu	ບຸJ	Serial dilution percent difference = 45.6%
		¢.	Baseline instability in prep blank
	Mg		Serial dilution percent difference = 14.9%
	Ni	ប	Baseline instability in cal blanks
	K .	J	Serial dilution percent difference = 42.7%
	Se	R	Analyte reported as potential false positive

May 21, 2003

INORGANIC DATA QUALIFIERS REPORT (continued)

Case Number: 31635 Project Number: 03-0474

<u>Sample No.</u>	Element	Flaq	Reason
6789	Al	J	Serial dilution percent difference = 13.4%
	Sb	R	Matrix spike recovery = 50.8%
•			Analyte reported as potential false positive
	Co ·	R	Analyte reported as potential false positive
	Cu	ŪJ	Serial dilution percent difference = 45.6%
			Baseline instability in prep blank
	Mg	J	Serial dilution percent difference = 14.9%
	Hg	U	Positive reported < lowest std on cal curve
	Ni	U	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
	Se,	R	Analyte reported as potential false positive
	Ag	R	Analyte reported as potential false positive
6790	Al	J	Serial dilution percent difference = 13.4%
•	Sb	R	Matrix spike recovery = 50.8%
	•		Analyte reported as potential false positive
	As	R	Analyte reported as potential false positive
	Cu	IJ	Serial dilution percent difference = 45.6%
			Baseline instability in prep blank
	Mg	· J	Serial dilution percent difference = 14.9%
	Ni ·	ับ	Baseline instability in cal blanks
	K	J	Serial dilution percent difference = 42.7%
	Ag	R	Analyte reported as potential false positive
	- 7	_	2 1 2 2 2 2 2 2
6791	Al	J	Serial dilution percent difference = 13.4%
	Sb	J	Matrix spike recovery = 50.8%
			Analyte reported as potential false positive
	Co	R	Analyte reported as potential false positive
	Cu .	UJ	Serial dilution percent difference = 45.6%
			Baseline instability in prep blank
	Mg	J	Serial dilution percent difference = 14.9%
	Ni .	. <b>U</b>	Baseline instability in cal blanks
	<b>K</b> .	<b>J</b>	Serial dilution percent difference = 42.7%
6792	Al	J	Serial dilution percent difference = 13.4%
	Sb	R	Matrix spike recovery = 50.8%
	The production of the state of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the section of the s		Analyte reported as potential false positive
• "	Cr	U	Baseline instability in cal blanks
	Co	R	Analyte reported as potential false positive
	Cu	J	Serial dilution percent difference = 45.6%
	Mg	J	Serial dilution percent difference = 14.9%
	Ni	Ū	Baseline instability in cal blanks
	. K	J	Serial dilution percent difference = 42.7%
	Ag	R	Analyte reported as potential false positive
	V	R	Analyte reported as potential false positive
6793	Al '	J	Serial dilution percent difference = 11.4%
	Cu	· U	Baseline instability in cal blanks
•	Fe	J	Matrix duplicate RPD = 53.5%
•	- Ni	ប	
	- N1	U	Baseline instability in cal blanks

May 21, 2003 INORGANIC DATA QUALIFIERS REPORT (continued)

Project Number: 03-0474

Sample No.	Element	Flag	Reason
6794	Al	J	Serial dilution percent difference = 11.4%
	Cu	U	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	Ū	Baseline instability in cal blanks
	Se	Ū	Baseline instability in cal blanks
		•	
6795	Al	Ĵ	Serial dilution percent difference = 11.4%
0,55	Cu	ับ	Baseline instability in cal blanks
•	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	Ŭ	Baseline instability in cal blanks
		Ŭ	baberine indeability in ear braine
6796	Al	J	Serial dilution percent difference = 11.4%
0750	Cu	Ū .	Baseline instability in cal blanks
	Fe .	J	Matrix duplicate RPD = 53.5%
	Ni Ni	Ŭ	Baseline instability in cal blanks
	Se	U	Baseline instability in cal blanks
•	Ag	R	Analyte reported as potential false positive
6707		~	Coming diluting manager difference 11 48
6797	Al	J	Serial dilution percent difference = 11.4%
	Cu	Ŭ	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Нg	ប 	Positive reported < lowest std on cal curve
	Ni	U	Baseline instability in cal blanks
•	Se	Ŭ	Baseline instability in cal blanks
• •	Na	J	% RSD > 20% for ICP multiple exposures
6700		· 🕶 ·	Camial dilubias massach differences 11 40
6798	A1	J.	Serial dilution percent difference = 11.4%
•	Cu	U -	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	ָּט	Baseline instability in cal blanks
,	Se	U	Baseline instability in cal blanks
6700	. 7 א	. т	Corial dilution pargent difference - 11 49
6799	Al	J	Serial dilution percent difference = .11.4%
•	Cu	Ŭ	Baseline instability in cal blanks
	ι Fe	J	Matrix duplicate RPD = 53.5%
•	Ni	Ŭ	Baseline instability in cal blanks
6800	Al	τ	Carial dilution pargent difference - 11 44
6800		J	Serial dilution percent difference = 11.4%
	Cu	Ŭ	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	Ŭ	Baseline instability in cal blanks
•	Se	J	% RSD > 20% for ICP multiple exposures
•	Na	J	% RSD > 20% for ICP multiple exposures
6901	וא	т	Corial dilution persont difference - 11 49
6801	Al	J	Serial dilution percent difference = 11.4%
	Cu	บ	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	ū	Baseline instability in cal blanks
	Ag	R	Analyte reported as potential false positive
•	Na	J	% RSD > 20% for ICP multiple exposures

May 21, 2003 INORGANIC DATA QUALIFIERS REPORT (continued)

Project Number: 03-0474

Sample No.	Element	Flag	Reason
6802	Al	J	Serial dilution percent difference = 11.4%
	Cu	Ū	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	U	Baseline instability in cal blanks
	Se	Ū	Baseline instability in cal blanks
6804	Al	J	Serial dilution percent difference = 11.4%
	Cu	U	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	U	Baseline instability in cal blanks
6805	Al	J	Serial dilution percent difference = 11.4%
	Cu	. U	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	Ŭ	Baseline instability in cal blanks
	Se	U	Baseline instability in cal blanks
6806	Al	J	Serial dilution percent difference = 11.4%
•	Cu	Ŭ	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
•	Ni	U	Baseline instability in cal blanks
6807	Al	J	Serial dilution percent difference = 11.4%
•	Cu	Ū	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	U	Baseline instability in cal blanks
	Ag	R	Analyte reported as potential false positive
•	Na	J	% RSD > 20% for ICP multiple exposures
6808	Al	J	Serial dilution percent difference = 11.4%
	Cu <sup>-</sup>	Ū	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	. U	Baseline instability in cal blanks
6809	Al	J	Serial dilution percent difference = 11.4%
	Cu	U	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
	Ni	Ŭ	Baseline instability in cal blanks
	Se ·	บ	Baseline instability in cal blanks
	Na	J	% RSD > 20% for ICP multiple exposures
6810	Al	J	Serial dilution percent difference = 11.4%
	Co	R	Analyte reported as potential false positive
	Cu	$\mathbf{U}^{\cdot}$	Baseline instability in cal blanks
	Fe	J	Matrix duplicate RPD = 53.5%
· · · · · · · · · · · · · · · · · · ·	Ni	U.	Baseline instability in cal blanks
	Na	J	% RSD > 20% for ICP multiple exposures

May 21, 2003 INORGANIC DATA QUALIFIERS REPORT (continued)

Case Number: 31635

Project Number: 03-0474
Site: Gulf States Creosoting, Flowood, MS

Sample No.	Element	Flag	Reason
6811	Al	J	PE sample recovery > warning limit
			PE sample recovery > action limit
,	Cu .	U	Baseline instability in blind blank
6812	Al	J	PE sample recovery > warning limit
			PE sample recovery > action limit
	Cu	Ū	Baseline instability in blind blank
6813	Al	J	PE sample recovery > warning limit
• •			PE sample recovery > action limit
6815	Al	IJ	PE sample recovery > warning limit
•			PE sample recovery > action limit
			Positive in blind blank
	Mg	U	Positives in cal, prep, and blind blanks
	K	Ü	Positive in blind blank
	Na	Ŭ	Positive in blind blank

Inorg Contractor: SENTIN

Sample 6753

Project: 03-0474 FY 2003

Produced by: Goddard, Denise Requestor:

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS06GW /

MD No: 1XR3

Project Leader: BSTRIGGO Beginning: 04/24/2003 13:45

Endina:

Media: GROUNDWATER

RESULTS	UNITS	ANALYTE
74	UG/L	Aluminum
2.4 U	UG/L	Antimony
3.5 U	UG/L	Arsenic
0.60 U	UG/L	Barium
0.10 U -	UG/L	Beryllium
0.20 U	UG/L	Cadmium
430 U	UG/L	Calcium
0.60 U	UG/L	Chromium
0.90 U	UG/L	Cobalt
0.67	UG/L	Copper
52 U	UG/L	Iron
2.0 U	UG/L	Lead
15 U	UG/L	Magnesium
0.50 U	UG/L	Manganese
0.10 U	UG/L	Total Mercury
1.5 U	UG/L	Nickei
100	UG/L	Potassium
2.4 U	UG/L	Selenium
0.70 U	UG/L	Silver
400	UG/L	Sodium
3.1 U	UG/L	Thallium
0.60 U	UG/L	Vanadium
11 U	UG/L	Zinc
NA	UG/L	Cyanide

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

# METALS SAMPLE ANALYSIS EPA - REGION IV SESD, ATHENS, GA

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Project: 03-0474 Sample 6754 FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:00 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS01SD / MD No: 1XR4 Org Contractor: LIBRTY D No: 1XR4 Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** 8200 J MG/KG Aluminum 1.2 UJ MG/KG Antimony 5.7 MG/KG Arsenic 89 MG/KG Barium 0.39 MG/KG Beryllium 0.10 U MG/KG Cadmium 1500 MG/KG Calcium 13 MG/KG Chromium 4.4 MG/KG Cobalt 10 U MG/KG Copper 15000 MG/KG Iron 12 J MG/KG Lead 1000 J MG/KG Magnesium 340 MG/KG Manganese 0.13 U MG/KG **Total Mercury** 6.6 U MG/KG Nickel 700 J MG/KG Potassium 1.9 R MG/KG Selenium 0.68 R MG/KG Silver 690 MG/KG Sodium 1.6 U MG/KG Thallium MG/KG Vanadium 24 36 MG/KG Zinc NA 'MG/KG Cyanide 61 % Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

WE I ALO SAMIFLE AMALTOID EPA - REGION IV SESD, ATHENS,

EPA - HEGION IV SESD, ATHENS, GA Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Project: 03-0474 Sample 6755 FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 08:50 Program: SF Case No: 31635 Ending: Id/Station: GS18SB / Inorg Contractor: SENTIN MD No: 1XR5 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XR5 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
7700 J	MG/KG	Aluminum
0.59 UJ	MG/KG	Antimony
4.1	MG/KG	Arsenic
26	MG/KG	Barium
0.31	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
190	MG/KG	Calcium
8.7	MG/KG	Chromium
3.4	MG/KG	Cobalt
3.9 U	MG/KG	Copper
12000	MG/KG	Iron
8.0 J	MG/KG	Lead
470 J	MG/KG	Magnesium
76	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
3.4 U	MG/KG	Nickel
220 J	MG/KG	Potassium
0.59 U	MG/KG	Selenium
0.32 R	MG/KG	Silver
540	MG/KG	Sodium
0.76 U	MG/KG	Thallium
18	MG/KG	Vanadium
16	MG/KG	Zinc
NA	MG/KG	Cyanide
18	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be ress than the reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### METALS SAMPLE ANALYSIS

### **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 06/05/2003 08:59

FY 2003 Sample 6756 Project: 03-0474

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS18SS /

Media: SURFACE SOIL (0" - 12")

MD No: 1XR6 D No: 1XR6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 08:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4900 J	MG/KG	Aluminum
1.0 UJ	MG/KG	Antimony
5.4	MG/KG	Arsenic
53	MG/KG	Barium
0.35	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
890	MG/KG	Calcium
9.9	MG/KG	Chromium
7.6	MG/KG	Cobalt
3.1 U	MG/KG	Copper
10000	MG/KG	iron
15 J	MG/KG	Lead
340 J	MG/KG	Magnesium
590	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
3.9 U	MG/KG	Nickel
180 J	MG/KG	Potassium
1.1	MG/KG	Selenium
0.29	MG/KG	Silver
310	MG/KG	Sodium
0.80 U	MG/KG	Thallium
20	MG/KG	Vanadium
34	MG/KG	Zinc
NA	MG/KG	Cyanide
22	·%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6757 FY 2003 Project: 03-0474

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS19SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XR7 D No: 1XR7

Inorg Contractor: SENTIN

**Org Contractor: LIBRTY** 

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:35

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
12000 J	MG/KG	Aluminum
0.79 UJ	MG/KG	Antimony
8.3	MG/KG	Arsenic
43	MG/KG	Barium
0.46	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
190	MG/KG	Calcium
12	MG/KG	Chromium
9.6	MG/KG	Cobalt
6.3 U	MG/KG	Copper
20000	MG/KG	Iron
12 J	MG/KG	Lead
680 J	MG/KG	Magnesium
600	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
5.4 U	MG/KG	Nickel
480 J 🔒	MG/KG	Potassium
1.7 J	MG/KG	Selenium
0.53	MG/KG	Silver
280	MG/KG	Sodium
0.78 U	MG/KG	Thallium
31	MG/KG	Vanadium
30	MG/KG	Zinc
NA	MG/KG	Cyanide
20	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

<b>METALS SA</b>	MPLE	ANAL	YSIS
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Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Project: 03-0474 Sample 6758 FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 09:30 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS19SS / MD No: 1XR8 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XR8 DATA REPORTED ON DRY WEIGHT BASIS

		and the second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second s
RESULTS	UNITS	ANALYTE
8500 J	MG/KG	Aluminum
0.89 UJ	MG/KG	Antimony
6.5	MG/KG	Arsenic
89	MG/KG	Barium
0.73	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
2000	MG/KG	Calcium
8.3	MG/KG	Chromium
13	MG/KG	Cobalt
3.8 U	MG/KG	Copper
12000	MG/KG	Iron
19 J	MG/KG	Lead
530 J	MG/KG	Magnesium
1900	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
6.7 U	MG/KG	Nickel
550 J	MG/KG	Potassium
1.2	MG/KG	Selenium
0.53 R	MG/KG	Silver
290	MG/KG	Sodium
0.73 U	MG/KG	Thallium
22	MG/KG	Vanadium
52	MG/KG	Zinc
NA	MG/KG	Cyanide
15	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 09:50 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS02SD / MD No: 1XR9 Org Contractor: LIBRTY Media: SEDIMENT D No: 1XR9 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4700 J	MG/KG	Aluminum
2.4 UJ	MG/KG	Antimony
. 1.8 R	MG/KG	Arsenic
77	MG/KG	Barium
0.56	MG/KG	Beryllium
0.09 U	MG/KG	Cadmium
880	MG/KG	Calcium
7.8	MG/KG	Chromium
4.7	MG/KG	Cobalt
6.6 U	MG/KG	Copper
6400	MG/KG	Iron
20 J	MG/KG	Lead
410 J	MG/KG	Magnesium
500	MG/KG	Manganese
0.11 U	MG/KG	Total Mercury
4.6 U	MG/KG	Nickel
410 J	MG/KG	Potassium
1.1 U	MG/KG	Selenium
0.45 R	MG/KG	Silver
580	MG/KG	Sodium
1.4 U	MG/KG	Thallium
13	MG/KG	Vanadium
33	MG/KG	Zinc
NA	MG/KG	Cyanide
56	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Sample FY 2003 Project: 03-0474 6760 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 10:05 Program: SF Case No: 31635 Ending: Id/Station: GS17SS / Inorg Contractor: SENTIN MD No: 1XS0 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XS0 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** 9000 J MG/KG Aluminum 0.83 UJ MG/KG **Antimony** MG/KG 5.5 Arsenic 120 MG/KG Barium 0.94 MG/KG Beryllium 0.07 R MG/KG Cadmium MG/KG Calcium 2300 8.0 MG/KG Chromium-12 MG/KG Cobalt 5.5 U MG/KG Copper MG/KG 12000 Iron 20 J MG/KG Lead 490 J MG/KG Magnesium 2500 MG/KG Manganese 0.08 U 'MG/KG **Total Mercury** 9.2 U MG/KG Nickel 620 J MG/KG Potassium 1.5 J· MG/KG Selenium 0.51 R MG/KG Silver 280 MG/KG Sodium 0.79 U MG/KG Thallium MG/KG 23 Vanadium 65 MG/KG Zinc NA MG/KG Cyanide % Moisture 21

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:20

Ending:

Program: SF ld/Station: GS17SB / Case No: 31635

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

MD No: 1XS1 D No: 1XS1

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
6700 J	MG/KG	Aluminum
1.1 UJ	MG/KG	Antimony
1.5	MG/KG	Arsenic
61	MG/KG	Barium
0.39	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
620	MG/KG	Calcium
7.0	MG/KG	Chromium
1.5 •	MG/KG	Cobalt
2.4 U	MG/KG	Copper
5900	MG/KG	Iron
9.1 J	MG/KG	Lead
310 J	MG/KG	Magnesium
57	MG/KG	Manganese
0.06 U	MG/KG	Total Mercur
2.5 U	MG/KG	Nickel
280 J	MG/KG	Potassium
0.60 U	MG/KG	Selenium
0.32 R	MG/KG	Silver
340	MG/KG	Sodium
0.78 U	MG/KG	Thallium
13	MG/KG	Vanadium
12	MG/KG	Zinc
NA	MG/KG	Cyanide
21	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474

**Metals Scan** 

Sample

Facility: Gulf States Creosoting

Media: SURFACE SOIL (0" - 12")

6762 FY 2003

Flowood, MS

Program: SF

Case No: 31635

ld/Station: GS20SS /

MD No: 1XS2 D No: 1XS2

Org Contractor: LIBRTY

Inorg Contractor: SENTIN

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:45

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
6700 J	MG/KG	Aluminum
0.60 UJ	MG/KG	Antimony
5.8	MG/KG	Arsenic
410	MG/KG	Barium
1.2	MG/KG	Beryllium
0.17	MG/KG	Cadmium
1700	MG/KG	Calcium
8.2	MG/KG	Chromium
17	MG/KG	Cobalt
0.95 U	MG/KG	Copper
12000	MG/KG	Iron
24 J	MG/KG	Lead
380 J	MG/KG	Magnesium
5500	MG/KG	Manganese
0.08 U	MG/KG	Total Mercury
8.0 U	MG/KG	Nickel
240 J	MG/KG	Potassium
1.3 J	MG/KG	Selenium
0.90 R	MG/KG	Silver
290	MG/KG	Sodium
1.5	MG/KG	Thallium
22	MG/KG	Vanadium
88 .	MG/KG	Zinc
NA	MG/KG	Cyanide
20	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates,

Produced by: Goddard, Denise Project: 03-0474 Sample 6763 FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 10:55 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS20SB / MD No: 1XS3 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XS3 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4700 J	MG/KG	Aluminum
0.57 UJ	MG/KG	Antimony
2.3	MG/KG	Arsenic
24	MG/KG	Barium
0.30	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
120	MG/KG	Calcium
4.7	MG/KG	Chromium
2.2	MG/KG	Cobalt
2.3 U	MG/KG	Copper
6900	MG/KG	Iron
6.0 J	MG/KG	Lead
210 J	MG/KG	Magnesium
140 .	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
1.7 U	MG/KG	Nickel
170 J	MG/KG	Potassium
0.57 U	MG/KG	Selenium
0.25 R	MG/KG	Silver
340	MG/KG	Sodium
0.74 U	MG/KG	Thallium
14	MG/KG	Vanadium
.11	MG/KG	Zinc
NA	MG/KG	Cyanide
16	%	% Moisture

# Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

METAL	S S	AMPL	E A	NAL	YSIS.
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Program: SF

Facility: Gulf States Creosoting

EPA - REGION IV SESD, ATHENS, GA

Production Date: 06/05/2003 08:59

Sample 6764 FY 2003 Project: 03-0474 Produced by: Goddard, Denise Requestor:

Metals Scan

Project: 03-0474

Requestor: RSTRIGGO

Flowood, MS Project Leader: BSTRIGGO
Beginning: 04/22/2003 10:45
Case No: 31635 Ending:

Id/Station: GS20SD / MD No: 1XS4 Inorg Contractor: SENTIN

Media: SEDIMENT D No: 1XS4 Org Contractor: LIBRTY DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
6100 J	MG/KG	Aluminum
0.59 UJ	MG/KG	Antimony
4.5	MG/KG	Arsenic
130	MG/KG	Barium
0.76 •	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
1400	MG/KG	Calcium
7.2	MG/KG	Chromium
6.8	MG/KG	Cobalt
4.3 U	MG/KG	Copper
8800	MG/KG	Iron
22 J	MG/KG	Lead
410 J	MG/KG	Magnesium
1800	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
6.1 U	MG/KG	Nickel
250 J	MG/KG	Potassium
1.1 R	MG/KG	Selenium
0.42	MG/KG	Silver
290	MG/KG	Sodium
0.77 U	MG/KG	Thallium
17	MG/KG	Vanadium
63	MG/KG	Zinc
NA	MG/KG	Cyanide
19	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474

**Metals Scan** 

Facility: Gulf States Creosoting

6765 FY 2003

Program: SF

Sample

Case No: 31635

Id/Station: GS03SD / Media: SEDIMENT

MD No: 1XS5

D No: 1XS5

Flowood, MS

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 10:30

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

DECLU TO	UNITS	ANALYTE
RESULTS		· · · · · · · · · · · · · · · · · · ·
16000·J	MG/KG	Aluminum
4.5 UJ	MG/KG	Antimony
7.9 R	MG/KG	Arsenic
300	MG/KG	Barium
1.3	MG/KG	Beryllium
0.25 U	MG/KG	Cadmium
2000	MG/KG	Calcium
- 24	MG/KG	Chromium
12	MG/KG	Cobalt
25 U	MG/KG	Copper
24000	MG/KG	Iron
41 J	MG/KG	Lead
1200 J	MG/KG	Magnesium
720	MG/KG	Manganese
0.31 U	MG/KG	Total Mercury
16 U	MG/KG	Nickel
1100 J	MG/KG	Potassium
4.2	MG/KG	Selenium
. 1.1 R	MG/KG	Silver
1500	MG/KG	Sodium
3.9 U	MG/KG	Thallium
41	MG/KG	Vanadium
130	MG/KG	Zinc
NA	MG/KG	Cyanide
84	% .	% Moisture
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Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate.

K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Sample 6766 FY 2003 Project: 03-0474 Requestor: **Metals Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 11:30 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS16SS / MD No: 1XS6 Org Contractor: LIBRTY D No: 1XS6 Media: SURFACE SOIL (0" - 12") DATA REPORTED ON DRY WEIGHT BASIS RESULTS LINITS ΔΝΔΙ ΥΤΕ

RESULIS	UNITS	ANALYIE
7700 J	MG/KG	Aluminum
0.70 UJ	MG/KG	Antimony
5.5 J	MG/KG	Arsenic
50	MG/KG	Barium
0.43	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
1200	MG/KG	Calcium
10	MG/KG	Chromium
4.0	MG/KG	Cobalt
5.3 U	MG/KG	Copper
12000	MG/KG	Iron
15 J	MG/KG	Lead
510 J	MG/KG	Magnesium
240	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
3.6 U	MG/KG	Nickel
490 J	MG/KG	Potassium
1.3	MG/KG	Selenium
0.39 R	MG/KG	Silver
270	MG/KG	Sodium
0.78 U	MG/KG	Thallium
22	MG/KG	Vanadium
33 <sup>.</sup>	MG/KG	Zinc
_ NA	MG/KG	Cyanide
21	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

METALS SAMPLE ANALYSIS

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise 6767 FY 2003 Project: 03-0474 Sample:

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

**D No: 1XS7** 

Id/Station: GS16SB /

MD No: 1XS7

Media: SUBSURFACE SOIL (> 12")

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:40

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
12000 J	MG/KG	Aluminum
0.62 UJ	MG/KG	Antimony
7.3	MG/KG	Arsenic
49	MG/KG	Barium
0.49	MG/KG	Beryllium
0.05 U .	MG/KG	Cadmium
390	MG/KG	Calcium
17	MG/KG	Chromium
2.1	MG/KG	Cobalt
7.7	MG/KG	Copper
23000	MG/KG	Iron ·
8.7 J	MG/KG	Lead
650 J	MG/KG	Magnesium
52	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
4.2 U	MG/KG	Nickel
470 J	MG/KG	Potassium
1.7 J	MG/KG	Selenium
0.45	MG/KG	Silver
290	MG/KG	Sodium
0.80 U	MG/KG	Thallium
33	MG/KG	Vanadium
25	MG/KG	Zinc
NA	MG/KG	Cyanide
23	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

<b>METALS</b>	SAMPLE	ANALYSIS
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Production Date: 06/05/2003 08:59

Sample **Metals Scan** 

Program: SF

6768 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 12:05

Ending:

Id/Station: GS21SS / Media: SURFACE SOIL (0" - 12")

Facility: Gulf States Creosoting

Flowood, MS Case No: 31635

MD No: 1XS8

Inorg Contractor: SENTIN

**D No: 1XS8** 

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
6800 J	MG/KG	Aluminum
0.96 UJ	MG/KG	Antimony
2.7	MG/KG	Arsenic
88	MG/KG	Barium
0.81	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
190	MG/KG	Calcium
5.6	MG/KG	Chromium
14	MG/KG	Cobalt
1.5 U	MG/KG	Copper
7300	MG/KG	Iron
13 J	MG/KG	Lead
270 J	MG/KG	Magnesium
2200	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
6.7 U	MG/KG	Nickel
250 J	MG/KG	Potassium
0.80	MG/KG	Selenium
0.52	MG/KG	Silver
. 300	MG/KG	Sodium
0.79 U	.MG/KG	Thallium
13	MG/KG	Vanadium
24	MG/KG	Zinc
NA	MG/KG	Cyanide
22	%	% Moisture

# Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6769 FY 2003

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS21SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XS9 D No: 1XS9

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
8600 J	MG/KG	Aluminum
0.60 UJ	MG/KG	Antimony
4.5	MG/KG	Arsenic
28	MG/KG	Barium
0.30	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
140	MG/KG	Calcium
14	MG/KG	Chromium
1.9	MG/KG	Cobalt
3.6	MG/KG	Copper
13000	MG/KG	Iron
6.8∙J	MG/KG	Lead
420 J	MG/KG	Magnesium
95	MG/KG	Manganese
0.06 U ,	-MG/KG	Total Mercury
3.6 U	MG/KG	Nickel
300 J	MG/KG	Potassium
0.95 R	MG/KG	Selenium
0.40	MG/KG	Silver
310	MG/KG	Sodium
0.78 U	MG/KG	Thallium
21	MG/KG	Vanadium
16	MG/KG	Zinc
NA	MG/KG	Cyanide
20	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Ident|fication of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Ident|fication of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/05/2003 08:59

Project: 03-0474 6770 FY 2003

**Metals Scan** 

Sample

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

Program: SF Id/Station: GS01GW /

MD No: 1XT0

D No: 1XT0

Inorg Contractor: SENTIN Org Contractor: A4

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 11:05

Ending:

Media: GROUNDWATER			
RESULTS	UNITS	ANALYTE	
320 UJ	UG/L	Aluminum	
2.4 U	UG/L	Antimony	
3.5 U	UG/L	Arsenic	
20.	UG/L	Barium	
0.11	UG/L	Beryllium	
0.20 U.	UG/L	Cadmium	
4600	UG/L	Calcium	
· 1.8 R	UG/L	Chromium	
1.2	UG/L	Cobalt	
1.5 U	UG/L	Copper	
1500	UG/L	Iron	
2.0 U	UG/L	Lead	
1500	UG/L	Magnesium	
23	UG/L	Manganese	
0.10 U	UG/L	Total Mercury	
1.5 U	UG/L	Nickel	
520	UG/L	Potassium	
2.4 U	UG/L	Selenium	
0.70 U	UG/L	Silver	
40000	UG/L	Sodium	
3.1 U	UG/L	Thallium	

Cyanide Analysis Not Requested

0.60 U

11 U

UG/L

UG/L

NA UG/L

Vanadium

Cyanide

Zinc

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates. R-Presence or absence of analyte can not be determined from data due to severe quality control problems. Data are rejected and considered unusable.

Project: 03-0474 6771 FY 2003 Sample

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 09:15

Ending:

Program: SF

Id/Station: GS01SB /

**Metals Scan** 

Flowood, MS

Case No: 31635

MD No: 1XT1

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

Facility: Gulf States Creosoting

Org Contractor: LIBRTY D No: 1XT1

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
7300 J	MG/KG	Aluminum
0.59 UJ	MG/KG	Antimony
6.7	MG/KG	Arsenic
62	MG/KG	Barium
0.31	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
850	MG/KG	Calcium
12	MG/KG	Chromium
3.3	MG/KG	Cobalt
6.4	MG/KG	Copper
13000	MG/KG	Iron
8.2 J	MG/KG	Lead
580 J	MG/KG	Magnesium
180	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
5.2 U	MG/KG	Nickel
320 J	MG/KG	Potassium
0.70	MG/KG	Selenium
0.38	MG/KG	Silver
290	MG/KG	Sodium
0.76 U	MG/KG	Thallium
22	MG/KG	Vanadium
18	MG/KG -	Zinc
NA	MG/KG	Cyanide
18 .	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6772 FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 09:00 Program: SF Case No: 31635 Ending: Id/Station: GS01SS / Inorg Contractor: SENTIN MD No: 1XT2 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT2 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4700 J	MG/KG	Aluminum
0.56 UJ	MG/KG	Antimony
6.3	MG/KG	Arsenic
72	MG/KG	Barium
0.33	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
1100	MG/KG	Calcium
8.4	MG/KG	Chromium
5.8	MG/KG	Cobalt
6.0	MG/KG	Copper
10000	MG/KG	Iron
11 J	MG/KG	Lead
520 J	MG/KG	Magnesium
630	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
4.8 Մ	MG/KG	Nickel
380 J	MG/KG	Potassium
1.3	MG/KG	Selenium
0.39 R	MG/KG	Silver
270	MG/KG	Sodium
0.73 U	MG/KG	Thallium
17	MG/KG	Vanadium
29	MG/KG	Zinc
NA	MG/KG	Cyanide
15	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6773 FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/22/2003 13:15 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS13SS / MD No: 1XT3 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT3 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4400 J	MG/KG	Aluminum
0.60 R	MG/KG	Antimony
6.3	MG/KG	Arsenic
58	MG/KG	Barium
0.54	MG/KG	Beryllium
0.16 R	MG/KG	Cadmium
1200	MG/KG	Calcium
7.8	MG/KG	Chromium
4.2	MG/KG	Cobalt
7.2 J	MG/KG	Copper
12000	MG/KG	Iron
33	MG/KG	Lead
280 J	MG/KG	Magnesium
940	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
5.2 U	MG/KG	Nickel
390 J	MG/KG	Potassium
1.2	MG/KG	Selenium
0.33 R	MG/KG	Silver
230	MG/KG	Sodium
0.69 U	MG/KG	Thallium
17	MG/KG	Vanadium
72	MG/KG	Zinc
NA	MG/KG	Cyanide
11 '	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyze analyzed in replicate. Reported value is "average" of replicates.

MET	ALS	SAM	DIF	ΔΝΔΙ	YSIS
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Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample Requestor:

**Metals Scan** 

Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:25 Program: SF Case No: 31635 Ending:

Inorg Contractor: SENTIN ld/Station: GS13SB / MD No: 1XT4

Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XT4 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
8400 J	MG/KG	Aluminum
0.58 R	MG/KG	Antimony
4.7	MG/KG	Arsenic
23	MG/KG	Barium
0.34	MG/KG	Beryllium
0.05 Ú	MG/KG	Cadmium
400	MG/KG	Calcium
. 11	MG/KG	Chromium
1.7	MG/KG	Cobalt
5.5 UJ	MG/KG	Copper
14000	MG/KG	Iron
5.1	MG/KG	Lead
470 J	MG/KG	Magnesium
82	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
4.3 U	MG/KG	Nickel
410 J	MG/KG	Potassium
0.57 U	MG/KG	Selenium
0.39	MG/KG	Silver
310	MG/KG	Sodium
0.73 U	MG/KG	Thallium
22	MG/KG	Vanadium
20	MG/KG	Zinc
NA	MG/KG	Cyanide
15 .	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474

FY 2003 6775 **Metals Scan** 

Requestor:

Facility: Gulf States Creosoting

Project Leader: BSTRIGGO Flowood, MS Beginning: 04/22/2003 11:35

Ending:

Program: SF Id/Station: GS04SD /

Sample

Case No: 31635

Inorg Contractor: SENTIN MD No: 1XT5

Org Contractor: LIBRTY D No: 1XT5 Media: SEDIMENT DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE					
9100 J	MG/KG	Aluminum					
2.1 UJ	MG/KG	Antimony	1				
5.5	MG/KG	Arsenic					
120	MG/KG	Barium					
1.2	MG/KG	Beryllium					
0.18 U	MG/KG	Cadmium					
2000	MG/KG	Calcium					
15	MG/KG	Chromium	ļ	-			
· 11	MG/KG	Cobalt					
12 UJ	MG/KG	Copper	:				
13000	MG/KG	Iron					
21	MG/KG	Lead	!				
1000 J	MG/KG	Magnesium					
280	MG/KG	Manganese					
0.22 U	MG/KG	Total Mercury	į				
15 U	MG/KG	Nickel	1				
811 J *	MG/KG	Potassium	1				
2.1 U	MG/KG	Selenium					
1.1 R	MG/KG	Silver					
1200	MG/KG	Sodium					
2.8 U	MG/KG	Thallium		•	•	-	
23	MG/KG	Vanadium					
84	MG/KG	Zinc					
NA	MG/KG	Cyanide Cyanide	1				
77	%	% Moisture	i				
			į.				

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

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Production Date: 06/05/2003 08:59

FY 2003 Sample Project: 03-0474 Produced by: Goddard, Denise

**Metals Scan** 

Requestor:

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:15

Program: SF

Ending:

Id/Station: GS05SD /

Case No: 31635 MD No: 1XT6

Inorg Contractor: SENTIN

Media: SEDIMENT D No: 1XT6 Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4900 J	MG/KG	Aluminum
2.1 R	MG/KG	Antimony
2.9 R	MG/KG	Arsenic
71	MG/KG	Barium
0.51	MG/KG	Beryllium
0.09 U	MG/KG	Cadmium
1200	MG/KG	Calcium
8.4	MG/KG	Chromium
5.5	MG/KG	Cobalt
5.8 UJ	MG/KG	Copper
8400	MG/KG	Iron
12	MG/KG	Lead
660 J	MG/KG	Magnesium
320	MG/KG	Manganese
0.11 U	MG/KG	Total Mercury
6.5 U	MG/KG	Nickel
480 J	MG/KG	Potassium
1.0 U	MG/KG	Selenium
0.71 R	MG/KG	Silver
600	MG/KG	Sodium
1.3 U	MG/KG	Thallium
13	MG/KG	Vanadium
39	MG/KG	Zinc
NA NA	MG/KG	Cyanide
54	%	% Moisture

# Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample FY 2003

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS07SD / Media: SEDIMENT

**Metals Scan** 

Flowood, MS

Case No: 31635

MD No: 1XT7 D No: 1XT7

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 12:45

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
140 J	MG/KG	Aluminum
0.61 UJ	MG/KG	Antimony
0.90 U	MG/KG	Arsenic
3.2	MG/KG	Barium
0.03 U	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
110 U	MG/KG	Calcium
0.61 U	MG/KG	Chromium
0.47	MG/KG	Cobalt
0.22 UJ	MG/KG	Copper
340	MG/KG	Iron
1.2 J	MG/KG	Lead
14 UJ	MG/KG	Magnesium
53	MG/KG	Manganese
0.06 U	MG/KG	Total Mercun
0.60 U	MG/KG	Nickel
52 J	MG/KG	Potassium
0.61 U	MG/KG	Selenium
0.18 U	MG/KG	Silver
330	MG/KG	Sodium
0.79 U	MG/KG	Thallium
0.57 R	MG/KG	Vanadium
2.9 U	MG/KG	Zinc
NA	MQ/KG	Cyanide
22	%	% Moisture
•		

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise

Requestor:

**Metals Scan** Project Leader: BSTRIGGO

Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 13:55 Program: SF Case No: 31635 Ending:

Inorg Contractor: SENTIN ld/Station: GS23SS / MD No: 1XT8 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XT8 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
7500 J	MG/KG	Aluminum
0.65 UJ	MG/KG	Antimony
5.0	MG/KG	Arsenic
83	MG/KG	Barium
0.50	MG/KG	Beryllium
0.10 R	MG/KG	Cadmium
960	MG/KG	Calcium
12	MG/KG	Chromium
6.8	MG/KG	Cobalt
. 12 J	MG/KG	Copper
13000	MG/KG	Iron
24	MG/KG	Lead
630 J	MG/KG	Magnesium
920	MG/KG	Manganese
0.08 U	MG/KG	Total Mercury
4.8 U	MG/KG	Nickel
480 J -	MG/KG	Potassium
1.1 R	MG/KG	Selenium
0.46 R	MG/KG	Silver
340	MG/KG	Sodium
0.83 U	MG/KG	Thallium
23	MG/KG	Vanadium
49	MG/KG	Zinc
NA	MG/KG	Cyanide
26	%	% Moisture

Sample

6778

FY 2003

### Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Requestor:

**Metals Scan** 

Facility: Gulf States Creosoting

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Program: SF

Sample

Id/Station: GS15SS /

Case No: 31635

MD No: 1XT9 D No: 1XT9

Org Contractor: LIBRTY

Inorg Contractor: SENTIN

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

Project Leader: BSTRIGGO

Beginning: 04/22/2003 13:40

<b>RESULTS</b>	UNITS	ANALYTE
8000 J	MG/KG	Aluminum
0.56 UJ	MG/KG	Antimony
6.9	MG/KG	Arsenic
72	MG/KG	Barium
0.71	MG/KG	Beryllium
0.10	MG/KG	Cadmium
2000	MG/KG	Calcium
11 .	MG/KG	Chromium
. 11	MG/KG	Cobalt
4.8 UJ	MG/KG	Copper
12000	MG/KG	Iron
26	MG/KG	Lead
430 J	MG/KG	Magnesium
1800	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
6.7 U	MG/KG	Nickel
580 J	MG/KG	Potassium
1.1 R	MG/KG	Selenium
0.50	MG/KG	Silver
.310	MG/KG	Sodium
0.73 U	MG/KG	Thallium
23	MG/KG	Vanadium
48	MG/KG	Zinc
NA	MG/KG.	Cyanide
15	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

6780 FY 2003

Project: 03-0474

Metals Scan

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS15SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XW0

D No: 1XW0

Flowood, MS

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:50

Endina:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
11000 J	MG/KG	Aluminum
0.61 UJ	MG/KG	Antimony
8.1	MG/KG	Arsenic
45	MG/KG	Barium
0.59	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
710	MG/KG	Calcium
21	MG/KG	Chromium
2.9	MG/KG	Cobalt
7.0 J	MG/KG	Copper
24000	MG/KG	Iron
9.2	MG/KG	Lead
640 J	MG/KG	Magnesium
410	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
5.8 U	MG/KG	Nickel
490 J	MG/KG	Potassium
2.3 J	MG/KG	Selenium
0.65	MG/KG	Silver
290	MG/KG	Sodium
0.79 U	MG/KG	Thallium
36	MG/KG	Vanadium
36	MG/KG	Zinc
NA	MG/KG	Cyanide
21	%	% Moisture

### Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 FY **2003** Sample

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ld/Station: GS02SS /

Case No: 31635

MD No: 1XW1

Media: SURFACE SOIL (0" - 12")

D No: 1XW1

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 13:13

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
7800 J	MG/KG	Aluminum
- 0.64 R	MG/KG	Antimony
3.3 J	MG/KG	Arsenic
120	MG/KG	Barium
0.94	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
690	MG/KG	Calcium
6.1	MG/KG	Chromium
17	MG/KG	Cobalt
2.0 J	MG/KG	Copper
8400	MG/KG	Iron
17	MG/KG	Lead
470 J	MG/KG	Magnesium
2800	MG/KG	Manganese
0.07 U	· MG/KG	Total Mercury
8.3 U	MG/KG	Nickel
240 J	MG/KG	Potassium
0.93 R <del>-</del>	MG/KG	Selenium
0.71 R	MG/KG	Silver
300	MG/KG	Sodium
0.81 U	MG/KG	Thallium
15	MG/KG	Vanadium
47	MG/KG	Zinc
NA	MG/KG	Cyanide
23	% .	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value. L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

**Metals Scan** 

Program: SF

6782

Facility: Gulf States Creosoting

FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 13:34

Ending:

Id/Station: GS02SB / Media: SUBSURFACE SOIL (> 12") Flowood, MS

Case No: 31635

MD No: 1XW2 D No: 1XW2

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
7800 J	MG/KG	Aluminum
0.59 UJ	MG/KG	Antimony
3.4 J	MG/KG	Arsenic
34	MG/KG	Barium
0.36	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
170	MG/KG	Calcium
9.5	MG/KG	Chromium
2.6 •	MG/KG	Cobalt
3.8 UJ	MG/KG	Copper
13000	MG/KG	Iron
6.6	MG/KG	Lead
480 J	MG/KG	Magnesium
200	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
4.1 U	MG/KG	Nickel
320 J	MG/KG	Potassium
1.4	MG/KG	Selenium
0.38 R	MG/KG	Silver
290	MG/KG	Sodium
0.77 U	MG/KG	Thallium
20 .	MG/KG	Vanadium
20	MG/KG	Zinc
NA	MG/KG	Cyanide
19	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:40

Ending:

Sample FY 2003 Project: 03-0474

**Metals Scan** 

Program: SF

Facility: Gulf States Creosoting

Id/Station: GS14SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635

MD No: 1XW3

Inorg Contractor: SENTIN

D No: 1XW3

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4800 J	MG/KG	Aluminum
0.69 J	MG/KG	Antimony
9.0	MG/KG	Arsenic
65	MG/KG	Barium
0.44	MG/KG	Beryllium
0.11 R	MG/KG	Cadmium
1800	MG/KG	Calcium
8.6	MG/KG	Chromium
7.3	MG/KG	Cobalt
5.7 UJ	MG/KG	Copper
11000	MG/KG	Iron
24	MG/KG	Lead
340 J	MG/KG	Magnesium
870	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
4.1 U	MG/KG	Nickel
200 J	MG/KG	Potassium
0.63 R	MG/KG	Selenium
0.37	MG/KG	Silver
300	MG/KG	Sodium
0.72 U	MG/KG	Thallium
21	MG/KG	Vanadium
44	MG/KG	Zinc
NA	MG/KG	~ j ~
14 .	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

MD No: 1XW4

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Sample 6784 FY **2003** Project: 03-0474 Requestor: **Metals Scan** Project Leader: BSTRIGGO

Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 14:50

Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS14SB /

Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XW4 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4600 J	MG/KG	Aluminum
0.56 UJ	MG/KG	Antimony
7.4	MG/KG	Arsenic
54	MG/KG	Barium
0.48	MĠ/KG	Beryllium
0.05 U	MG/KG	Cadmium
620	MG/KG	Calcium
7.1	MG/KG	Chromium <sup>*</sup>
12	MG/KG	Cobalt
1.5 UJ	MG/KG	Copper
14000	MG/KG	Iron
27	MG/KG	Lead
250 J	MG/KG	Magnesium
1000	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
3.3 U	MG/KG	Nickel
180 J	MG/KG	Potassium
0.56 U	MG/KG	Selenium
0.42	MG/KG	Silver
260	MG/KG	Sodium
0.73 U	MG/KG	Thallium
23	MG/KG	Vanadium
20	MG/KG	Zinc
NA	MG/KG	Cyanide
15	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Goddard, Denise Project: 03-0474 FY 2003 Sample **6785** Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 15:05 Program: SF Case No: 31635 Ending: Id/Station: GS10SS / Inorg Contractor: SENTIN MD No: 1XW5 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XW5

RESULTS	UNITS	ANALYTE
4700 J	MG/KG	Aluminum
1.0 R	MG/KG	Antimony
6.8	MG/KG	Arsenic
120	MG/KG	Barium
0.46	MG/KG .	Beryllium
0.34	MG/KG	Cadmium
2600	MG/KG.	Calcium
23	MG/KG	Chromium
4.8	MG/KG	Cobalt
19 J	MG/KG	Copper
14000	MG/KG	Iron
80	MG/KG	Lead
780 J	MG/KG	Magnesium
500	MG/KG	Manganese
0.06 ป	MG/KG	Total Mercury
8.5 U	MG/KG	Nickel
560 J	MG/KG	Potassium
0.98 R	MG/KG	Selenium
0.49	MG/KG	Silver
320	MG/KG	Sodium
0.70 U	MG/KG	Thallium
18	MG/KG	Vanadium
160	MG/KG	Zinc
NA	MG/KG	Cyanide
11	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6786 FY 2003 Project: 03-0474

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS10SB /

MD No: 1XW6

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

D No: 1XW6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 15:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	· UNITS	ANALYTE
8600 J	MG/KG	Aluminum
0.96 R	MG/KG	Antimony
13	MG/KG	Arsenic
430	MG/KG	Barium
1.3	MG/KG	Beryllium
0.87	MG/KG	Cadmium
9400	MG/KG	Calcium
19	MG/KG	Chromium
12	MG/KG	Cobalt ·
40 J	MG/KG	Copper
20000	MG/KG	Iron
71	MG/KG	Lead
2400 J	MG/KG	Magnesium
1800	MG/KG	Manganese
0.10 U	MG/KG	Total Mercury
20	MG/KG	Nickel
970 J	MG/KG	Potassium
1.2	MG/KG	Selenium
0.76	MG/KG	Silver
490	MG/KG	Sodium
0.76 U	MG/KG	Thallium
30	MG/KG	Vanadium
290	MG/KG	Zinc
NA	MG/KG	Cyanide
18	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample • 6787 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Metals Scan

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

MD No: 1XW7

Project Leader: BSTRIGGO Beginning: 04/22/2003 14:55

Ending:

Requestor:

Id/Station: GS02GW / Media: GROUNDWATER

Inorg Contractor: SENTIN Org Contractor: A4

D No: 1XW7

RESULTS	UNITS	ANALYTE
450 J	UG/L	Aluminum
2.4 U	UG/L	Antimony
3.5 U	UG/L	Arsenic
120	UG/L .	Barium
0.14	UG/L	Beryllium
0.20 U	UG/L	Cadmium
16000	UG/L	Calcium
0.92	UG/L	Chromium
1.2 R	UG/L	Cobalt
0.73 U	UG/L	Copper
1400	UG/L	Iron
2.0 U	UG/L	Lead
5500	UG/L	Magnesium
160	UG/L	Manganese
0.10 U	UG/L	Total Mercury
3.5	UG/L	Nickel
1500	UG/L	Potassium
2.4 U	UG/L	Selenium
0.70 U	.UG/L	Silver
120000	UG/L	Sodium
3.1 U	UG/L	Thallium
0.82 R	UG/L	Vanadium
13	UG/L	Zinc
NA	UG/L	Cyanide
•		

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable, reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6788 FY 2003 Project: 03-0474

**Metals Scan** 

Program: SF

Facility: Gulf States Creosoting

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:00

Ending:

Flowood, MS Case No: 31635

Id/Station: GS11SS / Media: SURFACE SOIL (0" - 12")

MD No: 1XW8 D No: 1XW8

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4400 J	MG/KG	Aluminum
0.51 UJ	MG/KG	Antimony
3.7	MG/KG	Arsenic
54	MG/KG	Barium
0.44	MG/KG	Beryllium
0.19	MG/KG	Cadmium
660	MG/KG	Calcium
12	MG/KG	Chromium
5.4	MG/KG	Cobalt
5.0 UJ	MG/KG	Copper
11000	MG/KG	Iron
290	MG/KG	Lead
260 J	MG/KG	Magnesium
880	MG/KG	Manganese
0.05 U	MG/KG	Total Mercury
4.9 U	MG/KG	Nickel
330 J	MG/KG	Potassium
0.77 R	MG/KG	Selenium
0.38	MG/KG	Silver
200	MG/KG	Sodium
0.66 U	MG/KG	Thallium
16	MG/KG	Vanadium
840	MG/KG	Zinc
NA	MG/KG	Cyanide
7	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample FY **2003** 

**Metals Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS11SB /

Media: SUBSURFACE SOIL (> 12")

Flowgood, MS

MD No: 1XW9

D No: 1XW9

Case No: 31635

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:10

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
9600 J	MG/KG	Aluminum
1.6 R	MG/KG	Antimony
6.6	MG/KG	Arsenic
41	MG/KG	Barium
0.49	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
320	MG/KG	Calcium
17	MG/KG	Chromium
3.2 R	MG/KG	Cobalt
5.5 UJ	MG/KG	Copper
20000	MG/KG	Iron
6.7	MG/KG	Lead
680 J	MG/KG	Magnesium
120	MG/KG	Manganese
0.10 U	MG/KG	Total Mercury
5.0 U	MG/KG	Nickel
430 J	MG/KG	Potassium
1.1 R	MG/KG	Selenium
1.1 R	MG/KG	Silver
350	MG/KG	Sodium
0.75 U	MG/KG	Thallium
33	MG/KG	Vanadium
44	MG/KG	Zinc
NA	MG/KG	Cyanide
18	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Project: 03-0474 Sample 6790 FY 2003

**Metals Scan** 

Facility: Gulf States Creosoting

Program: SF

ld/Station: GS09SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635

MD No: 1XX0

D No: 1XX0

Inorg Contractor: SENTIN Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2500 J	MG/KG	Aluminum
1.2 R	MG/KG	Antimony
0.95 R	MG/KG	Arsenic
25	MG/KG	Barium
0.15	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
340	MG/KG	Calcium
4.1	MG/KG	Chromium
0.95	MG/KG	Cobalt
0.87 UJ	MG/KG	Copper
2800	MG/KG	Iron
8.8	MG/KG	Lead
160 J	MG/KG	Magnesium
<b>28</b> .	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
1.5 U	MG/KG	Nickel
120 J	MG/KG	Potassium
0.60 U	MG/KG	Selenium
0.28 R 🗸	MG/KG	Silver
300	MG/KG	Sodium
0.77 U	MG/KG	Thallium
6.6	MG/KG	Vanadium
19	MG/KG	Zinc
NA	MG/KG	Cyanide
19	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

METALS SAMPLE ANALYSIS	METAL	S SAI	MPLE	ANAL	YSIS.
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Production Date: 06/05/2003 08:59

FY 2003 Project: 03-0474 Sample 6791

**Metals Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS09SB /

Media: SUBSURFACE SOIL (> 12")

Flowood, MS

Case No: 31635

MD No: 1XX1

D No: 1XX1

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:25

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
5800 J	MG/KG	Aluminum
0.60 UJ	MG/KG	Antimony
0.88 U	MG/KG	Arsenic
29	MG/KG	Barium
0.35	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
540	MG/KG	Całcium
5.3	MG/KG	Chromium
1.0 R	MG/KG	Cobalt
1.7 UJ •	MG/KG	Copper
3800	MG/KG	Iron
8.4	MG/KG	Lead
290 J	MG/KG	Magnesium
9.3	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
2.6 U	MG/KG	Nickel
190 J	MG/KG	Potassium
0.60 U	MG/KG	Selenium
0.18 U	MG/KG	Silver
440	MG/KG	Sodium
0.78 U	MG/KG	Thallium
8.4	MG/KG	Vanadium
12	MG/KG	Zinc .
NA	MG/KG	Cyanide
20	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

METALS SAMPLE ANALYSIS

6792 FY 2003

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 15:15

DATA REPORTED ON DRY WEIGHT BASIS

Ending:

Case No: 31635 Inorg Contractor: SENTIN MD No: 1XX2

D No: 1XX2

Flowood, MS

Project: 03-0474

Org Contractor: LIBRTY

**RESULTS UNITS ANALYTE** 

Facility: Gulf States Creosoting

Sample

**Metals Scan** 

Program: SF

Id/Station: GS06SD /

Media: SEDIMENT

450 J \_ MG/KG Aluminum 0.70 R MG/KG **Antimony** 0.88 U MG/KG Arsenic

6.8 MG/KG Barium 0.03 U Beryllium MG/KG 0.05 U MG/KG Cadmium

250 MG/KG Calcium 1.3 U MG/KG Chromium

0.61 R MG/KG Cobalt 0.15 UJ MG/KG Copper 950 MG/KG Iron 1.8 MG/KG Lead

MG/KG 64 J Magnesium 52 MG/KG Manganese 0.06 U MG/KG Total Mercury

0.76 U MG/KG Nickel 90 J MG/KG Potassium 0.60 U MG/KG Selenium 0.24 R MG/KG Silver 330

MG/KG Sodium 0.78 U MG/KG Thallium 1.6 R MG/KG Vanadium 3.6 MG/KG Zinc

NA MG/KG 20 %

Cyanide % Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences; | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

METALS SAMPLE ANALYSIS

**EPA - REGION IV SESD, ATHENS, GA** 

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Production Date: 06/05/2003 08:59

6793 FY 2003 Project: 03-0474 Sample

Facility: Gulf States Creosoting

Program: SF

Metals Scan

ld/Station: GS08SD / Media: SEDIMENT

Flowood, MS

Case No: 31635

MD No: 1XX3

D No: 1XX3

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/22/2003 15:50

Produced by: Goddard, Denise

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
510 J	MG/KG	Aluminum
0.65 U	MG/KG	Antimony
0.95 U	MG/KG	Arsenic
15	MG/KG	Barium
0.10	MG/KG	Beryllium
0.06 ป	MG/KG	Cadmium
180	MG/KG	Calcium
1.4	MG/KG	Chromium
1.0	MG/KG	Cobalt
1.2 U	MG/KG	Copper
2200 J	MG/KG	Iron
3.6	MG/KG	Lead
81	MG/KG	Magnesium
44	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
1.0 U	MG/KG	Nickel
56	MG/KG	Potassium
0.65 U	MG/KG	Selenium
0.19 U	MG/KG	Silver
68	MG/KG	Sodium
0.85 U	MG/KG	Thallium
2.4	MG/KG	Vanadium ·
4.6	MG/KG	Zinc
NA	MG/KG	Cyanide
27	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Project: 03-0474 Sample 6794 FY 2003 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/22/2003 16:42 Program: SF

Ending:

Id/Station: GS22SS / Inorg Contractor: SENTIN MD No: 1XX4 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XX4 DATA REPORTED ON DRY WEIGHT BASIS

Case No: 31635

RESULTS	UNITS	ANALYTE
4200 J	MG/KG	Aluminum
0.56 U	MG/KG	Antimony
4.0	MG/KG	Arsenic
42	MG/KG	Barium
0.43	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
280	MG/KG	Calcium
9.4	MG/KG	Chromium
5.5	MG/KG	Cobalt
2.5 U	MG/KG	Copper
9700 J	MG/KG	Iron
9.8	MG/KG	Lead
220	MG/KG	Magnesium
470	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
2.2 U	MG/KG	Nickel
170	MG/KG	Potassium
1.1 U	MG/KG	Selenium
0.20	MG/KG	Silver
44 U	MG/KG	Sodium
0.72 U	MG/KG	Thallium
21	MG/KG	Vanadium
19	MG/KG	Zinc
. NA	MG/KG	Cyanide
14	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not getected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

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Production Date: 06/05/2003 08:59

Sample 6795 FY **2003** Project: 03-0474 Produced by: Goddard, Denise

**Metals Scan** 

Requestor:

Facility: Gulf States Creosoting

Flowood, MS

Project Leader: BSTRIGGO

Program: SF

Beginning: 04/22/2003 16:40

Id/Station: GS06SS /

Case No: 31635

Ending:

Media: SURFACE SOIL (0" - 12")

MD No: 1XX5 D No: 1XX5

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
6700 J	MG/KG	Aluminum
0.65 U	MG/KG	Antimony
2.0	MG/KG	Arsenic
85	MG/KG	Barium
0.28	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
1200	MG/KG	Calcium
7.7	MG/KG	Chromium
1.4	MG/KG	Cobalt
4.9 U	MG/KG	Copper
5600 J	MG/KG	Iron
. 17	MG/KG	Lead
550	MG/KG	Magnesium
.170	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
2.8 U	MG/KG	Nickel
430	MG/KG	Potassium
0.65 U	MG/KG	Selenium
0.29	MG/KG	Silver
110	MG/KG	Sodium
0.84 U	MG/KG	Thallium
13	MG/KG	Vanadium
35	MG/KG	Zinc
NA	MG/KG	Cyanide
26	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

FY 2003 Project: 03-0474 6796

**Metals Scan** 

Sample

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

Id/Station: GS06SB /

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XX6 D No: 1XX6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:50

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
7700 J	MG/KG	Aluminum
0.65 U	MG/KG	Antimony
2.7	MG/KĢ	Arsenic
48	MG/KG	Barium
0.24	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
990	MG/KG	Calcium
7.5	MG/KG	Chromium
1.2	MG/KG	Cobalt
3.4 U	MG/KG	Copper
5800 J	MG/KG	Iron
11	MG/KG	Lead
580	MG/KG	Magnesium
18	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
2.7 U	MG/KG	Nickel
240	MG/KG	Potassium
0.96 U	MG/KG	Selenium
0.28 R	MG/KG	Silver
250 .	MG/KG	Sodium
0.84 U	MG/KG	Thallium
. 17	MG/KG	Vanadium
12	MG/KG	Zinc
NA	MG/KG	Cyanide
27	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

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Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Sample 6797 FY 2003 Project: 03-0474

**Metals Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS24SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635

MD No: 1XX7

D No: 1XX7

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Requestor:

Project Leader: BSTRIGGO Beginning: 04/22/2003 16:35

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4900 J	MG/KG	Aluminum
0.91 U	MG/KG	Antimony
6.5	MG/KG	Arsenic
120	MG/KG	Barium
0.61	MG/KG	Beryllium
0.16	MG/KG	Cadmium
1600	MG/KG	Calcium
12	MG/KG	Chromium
6.8	MG/KG	Cobalt
8.4 U	MG/KG	Copper
17000 J .	MG/KG	Iron
28	MG/KG	Lead
450	MG/KG	Magnesium
890	MG/KG	Manganese
0.09 U	MG/KG	Total Mercury
5.3 U	MG/KG	Nickel
370	MG/KG	Potassium
1.5 U	MG/KG	Selenium
0.47	MG/KG	Silver
86 J	MG/KG	Sodium
1.2 U	MG/KG	Thallium
26	MG/KG	Vanadium
<sub>.</sub> 89	MG/KG	Zinc
NA	MG/KG	Cyanide
47	% .	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

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Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise Sample. 6798 FY 2003 Project: 03-0474 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 07:40 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS12SS / MD No: 1XX9 Media: SURFACE SOIL (0" - 12") Org Contractor: LIBRTY D No: 1XX9 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
5500 J	MG/KG	Aluminum
0.58 U 🕝	MG/KG	Antimony
4.4	MG/KG	Arsenic
83	MG/KG	Barium
0.46	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
760	MG/KG	Calcium
8.9	MG/KG	Chromium
7.3	MG/KG	Cobalt
3.4 U	MG/KG	Copper
11000 J	MG/KG	Iron
25	MG/KG	Lead
380	MG/KG	Magnesium
990	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
3.2 U	MG/KG	Nickel
280	MG/KG	Potassium
1.1 U	MG/KG	Selenium
0.43	MG/KG	Silver
91	MG/KG	Sodium
0.75 U	MG/KG	Thallium
22	MG/KG	Vanadium
28	MG/KG	Zinc
NA	MG/KG	Cyanide
18	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable, reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences! | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise . 6799 FY 2003 Project: 03-0474 Sample Requestor: **Metals Scan** Project Leader: BSTRIGGO Flowood, MS Facility: Gulf States Creosoting Beginning: 04/23/2003 07:50 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS12SB / MD No: 1XY0 Org Contractor: LIBRTY Media: SUBSURFACE SOIL (> 12") D No: 1XY0 DATA REPORTED ON DRY WEIGHT BASIS **RESULTS UNITS ANALYTE** 3700 J MG/KG Aluminum 0.58 U MG/KG Antimony MG/KG 4.6 Arsenic 19 MG/KG Barium 0.30 MG/KG Beryllium 0.05 U MG/KG Cadmium 160 MG/KG Calcium MG/KG Chromium 8.5 1.4 MG/KG Cobalt 1.8 U MG/KG Copper 13000 J MG/KG Iron 6.7 MG/KG Lead 170 MG/KG Magnesium 55 MG/KG Manganese 0.06 U MG/KG Total Mercury 1.4 U MG/KG Nickel 170 MG/KG Potassium MG/KG Selenium 1.4 0.47 MG/KG Silver 96 MG/KG Sodium 0.75 U MG/KG Thallium MG/KG 26 Vanadium 8.4 MG/KG Zinc NA MG/KG Cyanide % Moisture

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

Cyanide Analysis Not Requested

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

METALS SAMPLE ANALYSIS

Production Date: 06/05/2003 08:59

Produced by: Goddard, Denise FY 2003 Project: 03-0474 Sample 6800 Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 07:40 Program: SF Case No: 31635 Ending: Inorg Contractor: SENTIN Id/Station: GS03SS / MD No: 1XY1 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XY1 DATA REPORTED ON DRY WEIGHT BASIS RESULTS UNITS **ANALYTE** 8900 J MG/KG Aluminum 0.64 U MG/KG **Antimony** 11 MG/KG Arsenic 67 MG/KG Barium 0.94 MG/KG Beryllium 0.05 U MG/KG Cadmium 790 Calcium MG/KG 20 MG/KG Chromium 8.1 MG/KG Cobalt 6.5 U MG/KG Copper 27000 J MG/KG Iron 120 MG/KG Lead 680 MG/KG Magnesium 550 MG/KG Manganese 0.07 U MG/KG **Total Mercury** 5.1 U MG/KG Nickel

Cyanide Analysis Not Requested

350

0.63

1.9 J

82 J

0.83 U

46

34

25

NA

MG/KG

MG/KG

MG/KG

MG/KG

MG/KG

MG/KG

MG/KG

MG/KG

Potassium

Selenium

Silver

Zinc

Sodium

Thallium

Cvanide

% Moisture

Vanadium

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6801 FY 2003 Project: 03-0474

**Metals Scan** 

Facility: Gulf States Creosoting

Flowood, MS

Program: SF

ng nowood, n

Id/Station: GS03SB /

MD No. 1770

Media: SUBSURFACE SOIL (> 12")

Case No: 31635

MD No: 1XY2 D No: 1XY2 Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 07:55

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
8600 J	MG/KG	Aluminum
0.67 U	MG/KG	Antimony
9.4	MG/KG	Arsenic
63	MG/KG	Barium
0.64	MG/KG	Beryllium
0.06 U	MG/KG	Cadmium
1400	MG/KG	Calcium
14	MG/KG	Chromium
4.3	MG/KG	Cobalt
6.0 U	MG/KG	Copper
22000 J	MG/KG	Iron
15	MG/KG	Lead -
770	MG/KG	Magnesium
200	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
5.1 U	MG/KG	Nickel
360	MG/KG	Potassium
1.6	MG/KG	Selenium
0.49 R	MG/KG	Silver
. 98 J	MG/KG	Sodium
0.87 U	MG/KG	Thallium
33	MG/KG	Vanadium
23	MG/KG	Zinc
NA	MG/KG	Cyanide
28	. %	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit. N-Presumptive evidence analyte is present; analyte reported as tentative identification. NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### **METALS SAMPLE ANALYSIS**

## **EPA - REGION IV SESD, ATHENS, GA**

Production Date: 06/05/2003 08:59

Sample FY 2003 Project: 03-0474 6802

**Metals Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS08SS /

Media: SURFACE SOIL (0" - 12")

Flowood, MS

Case No: 31635

MD No: 1XY3

D No: 1XY3

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:10

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
1800 J	MG/KG	Aluminum
0.50 U	MG/KG	Antimony
3.4	MG/KG	Arsenic
21	MG/KG	Barium
0.19	MG/KG	Beryllium
0.09	MG/KG	Cadmium
800	MG/KG	_ Calcium _
4.0	MG/KG	Chromium
1.5	MG/KG	Cobalt
2.9 U	MG/KG	Copper
4700 J	MG/KG	Iron
12 .	MG/KG	Lead
180	MG/KG	Magnesium
230	MG/KG	Manganese
0.05 U	MG/KG	Total Mercury
1.7 U	MG/KG	Nickel
150	MG/KG	Potassium
0.68 U	MG/KG	Selenium
0.15 U	MG/KG	Silver
39 U	MG/KG	Sodium
0.65 U	MG/KG	Thallium
7.1 °	MG/KG	Vanadium
26	MG/KG	Zinc
NA	MG/KG	Cyanide
5	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/05/2003 08:59

Sample **Metals Scan** 

6804

FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/23/2003 08:20

Ending:

Program: SF

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

Id/Station: GS08SB /

MD No: 1XY4

D No: 1XY4

Inorg Contractor: SENTIN

Media: SUBSURFACE SOIL (> 12")

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
8100 J	MG/KG	Aluminum
0.58 U	MG/KG	Antimony
3.5	MG/KG	Arsenic
45	MG/KG	Barium
0.59	MG/KG	Beryllium -
0.05 U	MG/KG	Cadmium
450	MG/KG	Calcium
9.2	MG/KG	Chromium
9.4	MG/KG	Cobalt
3.9 U	MG/KG	Copper
12000 J	MG/KG	Iron
5.7	MG/KG	Lead
370	MG/KG	Magnesium
1000	MG/KG	Manganese
0:06 U	MG/KG	Total Mercury
5.3 U	MG/KG	Nickel
350	MG/KG	Potassium
0.58 U	MG/KG	Selenium
0.32	MG/KG	Silver
46	MG/KG .	Sodium
0.75 U	MG/KG	Thallium
22	MG/KG	Vanadium
25	MG/KG	Zinc
NA	MG/KG	Cyanide
18	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Produced by: Goddard, Denise Project: 03-0474 Sample 6805 **EY 2003** Requestor: **Metals Scan** Project Leader: BSTRIGGO Facility: Gulf States Creosoting Flowood, MS Beginning: 04/23/2003 08:35 Program: SF Case No: 31635 Ending: Id/Station: GS07SS / Inorg Contractor: SENTIN MD No: 1XY5 Org Contractor: LIBRTY Media: SURFACE SOIL (0" - 12") D No: 1XY5 DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
2800 J	MG/KG	Aluminum
0.58 U	MG/KG	Antimony
3.6	MG/KG	Arsenic
34	MG/KG	Barium
0.20	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
810	MG/KG	Calcium
4.5	MG/KG	Chromium
2.0	MG/KG	Cobalt
3.5 U	MG/KG	Copper
5400 J	MG/KG	Iron
11	MG/KG	Lead
250	MG/KG	Magnesium
180	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
1.9 U	MG/KG	Nickel
170	MG/KG	Potassium
0.79 U	MG/KG	Selenium
0.17 U	MG/KG	Silver
63	MG/KG	Sodium
0.75 U	MG/KG	Thallium
	MG/KG	<u>V</u> anadium
24	MG/KG	Zinc
NA	MG/KG	Cyanide
17	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

#### **METALS SAMPLE ANALYSIS**

Production Date: 06/05/2003 08:59

Sample 6806

FY 2003 Project: 03-0474

Facility: Gulf States Creosoting

**Metals Scan** 

Program: SF Id/Station: GS07SB /

Media: SUBSURFACE SOIL (> 12")

Flowood, MS

Case No: 31635

MD No: 1XY6 .

D No: 1XY6

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:45

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
9400 J	MG/KG	Aluminum
0.65 U	MG/KG	Antimony
2.7	MG/KG	Arsenic
50	MG/KG	Barium
0.35	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
670	MG/KG	Calcium
8.0	MG/KG	Chromium
1.6	MĢ/KG	Cobalt
5.7 U	MG/KG	Copper
8200 J	MG/KG	Iron
11	MG/KG	Lead ·
550	MG/KG	Magnesium
19	:MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
3.7 U	MG/KG	Nickel
260	MG/KG	Potassium
0.65 U	MG/KG	Selenium
0.19 U	MG/KG	Silver
220	MG/KG	Sodium
0.84 U	MG/KG	Thallium
18 *	MG/KG	Vanadium
15	MG/KG	Zinc
NA	MG/KG	Cyanide
26	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

**Metals Scan** 

Program: SF

FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 08:30

Ending:

Facility: Gulf States Creosoting Flowood, MS

ld/Station: GS04SS /

Case No: 31635

Media: SURFACE SOIL (0" - 12")

MD No: 1XY7 D No: 1XY7

Inorg Contractor: SENTIN Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
4300 J	MG/KG	Aluminum
0.63 U	MG/KG	Antimony
2.0	MG/KG	Arsenic
60	MG/KG	Barium
0.37	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
690	MG/KG	Calcium
6.0	MG/KG	Chromium
1.7	MG/KG	Cobalt
3.9 U	MG/KG	Copper
4500 J	MG/KG	Iron
18	MG/KG	Lead
280	MG/KG	Magnesium
140	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
2.6 U	MG/KG	Nickel
200	MG/KG -	Potassium
0.63 U	MG/KG	Selenium
0.26 R	MG/KG	Silver
120 J	MG/KG	Sodium
0.81 U	MG/KG	Thallium
11 '	MG/KG	Vanadium
27	MG/KG	Zinc
_ NA	MG/KG	Cyanide
23	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Facility: Gulf States Creosoting

Media: SUBSURFACE SOIL (> 12")

Sample

**Metals Scan** 

Program: SF

Id/Station: GS04SB /

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/05/2003 08:59

6808 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO

Beginning: 04/23/2003 08:40

Ending:

Flowood, MS

Case No: 31635 MD No: 1XY8

D No: 1XY8

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
6600 J	MG/KG	Aluminum
0.66 U	MG/KG	Antimony
1.9	MG/KG	Arsenic
57 <b>·</b>	MG/KG	Barium
0.41	MG/KG	Beryllium
0.06 U	MG/KG	Cadmium
670 .	MG/KG	Calcium
6.6	MG/KG	Chromium
1.4	MG/KG	Cobalt
4.0 U	MG/KG	Copper
7500 J	MG/KG	Iron .
9.2	MG/KG	Lead
370	MG/KG	Magnesium
60	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
3.0 U	MG/KG	Nickel
190	MG/KG	Potassium
0.66 U	MG/KG	Selenium
0.19 U	MG/KG	Silver
300	MG/KG	Sodium
0.85 U	MG/KG	Thallium
14	MG/KG	Vanadium
15	MG/KG	Zinc
NA	MG/KG	Cyanide
·27	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**METALS SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Production Date: 06/05/2003 08:59

FY **2003** Project: 03-0474 Sample 6809

**Metals Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS05SS / Media: SURFACE SOIL (0" - 12") Flowood, MS

Case No: 31635

MD No: 1XY9

D No: 1XY9

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:15

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
3100 J	MG/KG	Aluminum
0.57 U	MG/KG	Antimony
13	MG/KG	Arsenic
37	MG/KG	Barium
0.30	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
500	MG/KG	Calcium
7.5	MG/KG	Chromium
6.6	MG/KG	Cobalt
3.0 U	MG/KG	Copper
15000 J	MG/KG	Iron
19	MG/KG	Lead
180	MG/KG	Magnesium
610	MG/KG	Manganese
0.06 U	MG/KG	Total Mercury
2.0 U	MG/KG	Nickel
120	MG/KG	Potassium
0.87 U	MG/KG	Selenium
0.43	MG/KG	Silver
. 58 J	MG/KG	Sodium
0.73 U	MG/KG	Thallium
18 ,	MG/KG	Vanadium
23	MG/KG	Zinc
NA	MG/KG	Cyanide
15	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value,

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/05/2003 08:59

Project: 03-0474 Sample 6810 FY 2003

**Metals Scan** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS05SB /

Media: SUBSURFACE SOIL (> 12")

Flowood, MS

Case No: 31635

MD No: 1XZ0

D No: 1XZ0

Inorg Contractor: SENTIN

Org Contractor: LIBRTY

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:25

Ending:

DATA REPORTED ON DRY WEIGHT BASIS

RESULTS	UNITS	ANALYTE
5300 J	MG/KG	Aluminum
0.62 U	MG/KG	Antimony
0.91 U	MG/KG	Arsenic
59	MG/KG	Barium
0.27	MG/KG	Beryllium
0.05 U	MG/KG	Cadmium
580	MG/KG	Calcium
4.9	MG/KG	Chromium
0.78 R	MG/KG	Cobalt
2.1 U	MG/KG	Copper
2300 J	MG/KG	Iron
6.7	MG/KG	Lead
220 .	MG/KG	Magnesium
16	MG/KG	Manganese
0.07 U	MG/KG	Total Mercury
1.7 U	MG/KG	Nickel
170	MG/KG	Potassium
0.62 U	MG/KG	Selenium
0.18 U	MG/KG	Silver
190 J	MG/KG	Sodium
0.81 U	MG/KG	Thallium
6.7	MG/KG	Vanadium
7.3	MG/KG	Zinc
NA	MG/KG	Cyanide
23	%	% Moisture

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample

FY 2003 6811

Project: 03-0474

Produced by: Goddard, Denise

**Metals Scan** 

Program: SF

Facility: Gulf States Creosoting

Flowood, MS

Case No: 31635

Project Leader: BSTRIGGO Beginning: 04/23/2003 09:55

Id/Station: GS03GW /

MD No: 1XZ1

Ending:

Requestor:

Media: GROUNDWATER

D No: 1XZ1

Inorg Contractor: SENTIN

Org Contractor: A4

RESULTS	UNITS	ANALYTE
850 J	UG/L	Aluminum
2.4 U	UG/L	Antimony
3.5 U	UG/L	Arsenic
29 .	UG/L	Barium
0.11	UG/L ,	Beryllium
0.20 U	UG/L	Cadmium
3600	UG/L	Calcium
1.8	UG/L	Chromium
0.90 U	UG/L	Cobalt
1.1 U	UG/L	Copper
2100	UG/L	iron
2.0 U	UG/L	Lead
940	UG/L	Magnesium
110	UG/L	Manganese
0.10 U	UG/L	Total Mercury
1.5 U	UG/L	Nickel
870	UG/L	Potassium
2.4 U	UG/L	Selenium
0.70 U	UG/L	Silver
	UG/L	Sodium
3.1 U	UG/L	Thallium
2.0	UG/L	Vanadium
18	UG/L	Zinc
NA	UG/L	Cyanide

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**METALS SAMPLE ANALYSIS** 

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/05/2003 08:59

Sample

6812 F

FY 2003 Project: 03-0474

Produced by: Goddard, Denise

Requestor:

Project Leader: BSTRIGGO Beginning: 04/23/2003 13:20

Ending:

Metals Scan

Program: -SF

Facility: Gulf States Creosoting

Flowgod, MS

WDOG, IVIS

Case No: 31635 MD No: 1XZ3

Inorg Contractor: SENTIN

Id/Station: GS04GW / Media: GROUNDWATER

D No: 1XZ3

Org Contractor: A4

RESULTS	UNITS	ANALYTE
1700 J	UG/L	Aluminum
2.4 U	UG/L	Antimony
3.5 U	UG/L	Arsenic
26	UG/L	Barium
0.10 U	UG/L	Beryllium
0.20 U	UG/L	Cadmium
3500	UG/L	Calcium
1.6	UG/L	Chromium
0.90 U	UG/L	Cobalt
0.78 U	UG/L	Copper
1900	UG/L ·	lron
2.0 U	UG/L	Lead
930	UG/L	Magnesium
30	UG/L	Manganese
0.10 U	UG/L	Total Mercury
3.3	UG/L	Nickel
1800	UG/L	Potassium ,
2.4 U	UG/L	Selenium
0.70 U	UG/L	Silver
5000	UG/L	Sodium
3.1 U	UG/L	Thallium
1.9	UG/L	Vanadium
11 U ·	UG/L	Zinc
NA	UG/L	Cyanide

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Production Date: 06/05/2003 08:59

Sample

FY **2003** 

Project: 03-0474

**Metals Scan** 

Facility. Gulf States Creosoting

Flowood, MS

Program: SF

Case No: 31635

Id/Station: GS05GW /

MD No: 1XZ4

Inorg Contractor: SENTIN

Media: GROUNDWATER

D No: 1XZ4

Org Contractor: A4

Project Leader: BSTRIGGO Beginning: 04/23/2003 11:45

Produced by: Goddard, Denise

Ending:

Requestor:

RESULTS	UNITS	ANALYTE
370 J	UG/L	Aluminum
2.4 U	UG/L	Antimony
3.5 U	UG/L	Arsenic
32	UG/L	Barium
0.90	UG/L	Beryllium
0.20 U	UG/L	Cadmium
15000	UG/L	Calcium
0.60 U	UG/L	Chromium .
10	UG/L	Cobalt
0.60 U	UG/L	Copper
12000	UG/L	Iron
2.0 U	UG/L	Lead
8200	UG/L	Magnesium
320	UG/L	Manganese
0.10 U	UG/L	Total Mercury
20	UG/L	Nickel
1900	UG/L	Potassium
2.4 U	UG/L	Selenium
0.70 U	UG/L	Silver
86000	UG/L	Sodium
3.1 U	UG/L	Thallium
0.60 U	UG/L	Vanadium
42	UG/L	Zinc
· NA	UG/L	Cyanide

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

Sample 6815 FY 2003

Project: 03-0474

Produced by: Goddard, Denise

Requestor:

**Metals Scan** 

Flowood, MS Facility: Gulf States Creosoting

Program: SF

Project Leader: BSTRIGGO

Case No: 31635

Beginning: 04/24/2003 14:00

Id/Station: GS01PB /

MD No: 1XZ5

Ending:

Media: PRESERVATIVE BLANK

Inorg Contractor: SENTIN

UNITS	ANALYTE
UG/L	Aluminum
UG/L	Antimony
UG/L	Arsenic
UG/L	Barium
UG/L	Beryllium
UG/L	Cadmium
UG/L	Calcium
UG/L	Chromium
UG/L	Cobalt
UG/L	Copper
UG/L	lron .
UG/L	Lead
UG/L	Magnesium
UG/L	Manganese
UG/L	Total Mercury
UG/L	Nickel
UG/L	Potassium
UG/L	Selenium
UG/L	Silver
	Sodium
	Thallium
UG/L	Vanadium
UG/L	Zinc
UG/L	Cyanide
	UG/L UG/L UG/L UG/L UG/L UG/L UG/L UG/L

Cyanide Analysis Not Requested

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

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NA-Not Analyzed. | NAI-Not Analyzed due to Interferences: | A-Analyte analyzed in replicate. Reported value is "average" of replicates.



## UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

#### Region 4

Science and Ecosystem Support Division 980 College Station Road Athens, Georgia 30605-2720

#### **MEMORANDUM**

Date: 06/09/2003

Subject: Results of CLASSICALS/NUTRIENTS Sample Analysis

03-0558 Gulf States Creosoting

Flowood, MS

From: Howes, Roberta

To: Striggow, Brian

Thru: Scifres, Jenny

Chief, Inorganic Chemistry Section

**Analytical Support Branch** 

Attached are the results of analysis of samples collected as part of the subject project. If you have any questions, please contact me.

## Sample Disposal Policy:

According to our records this project is not part of a criminal investigation. Because of our limited space for long term sample storage, we must perform disposals on a routine basis.

Therefore, please take note that within 90 days of the date of this memo, the original samples and all extracts associated with the samples will be disposed of as required by all applicable and appropriate statutes.

These samples may be held in custody for longer than 90 days only by contacting our sample coordinator, Debbie Colquitt, by e-mail at Colquitt.Debbie@epa.gov.

#### ATTACHMENT.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/09/2003 15:36

Sample

6693 FY 2003

Project: 03-0558

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Program: SF

2.0

Id/Station: GS-01-SD / Media: SEDIMENT

Flowood, MS

Ending:

samples placed in custody room cooler 04/25/03 @ 1440 by Bria

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Howes, Roberta

Project Leader: BSTRIGGO

Beginning: 04/22/2003 09:00

Requestor: Brian Farrier

**RESULTS UNITS** 

%

% Total Organic Carbon

**ANALYTE** 

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/09/2003 15:36

Sample

6694

FY 2003

Project: 03-0558

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS-02-SD / Media: SEDIMENT

Flowood, MS

Ending:

samples placed in custody room cooler 04/25/03 @ 1440 by Bria

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Howes, Roberta

Project Leader: BSTRIGGO

Beginning: 04/24/2003 09:50

Requestor: Brian Farrier

RESULTS UNITS

2.1 %

% Total Organic Carbon

**ANALYTE** 

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/09/2003 15:36

Sample

6695 FY 2003

Project: 03-0558

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Program: SF

9.5

ld/Station: GS-03-SD / Media: SEDIMENT

Flowood, MS

Produced by: Howes, Roberta Requestor: Brian Farrier Project Leader: BSTRIGGO Beginning: 04/22/2003 10:30

Ending:

samples placed in custody room cooler 04/25/03 @ 1440 by Bria

DATA REPORTED ON DRY WEIGHT BASIS

**RESULTS UNITS** 

% Total Organic Carbon

ANALYTE

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/09/2003 15:36

Sample

6696 FY 2003

Project: 03-0558

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Program: SF

id/Station: GS-05-SD / Media: SEDIMENT

Flowood, MS

Beginning: 04/22/2003 12:15 Ending:

samples placed in custody room cooler 04/25/03 @ 1440 by Bria

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Howes, Roberta

Requestor: Brian Farrier

Project Leader: BSTRIGGO

**RESULTS UNITS** 

1.8 % **ANALYTE** % Total Organic Carbon

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification-of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/09/2003 15:36

Sample

6697

FY 2003

Project: 03-0558

**SPECIFIED TESTS** 

Facility: Gulf States Creosoting

Program: SF

Id/Station: GS-06-SD / Media: SEDIMENT Flowood, MS

Ending:

samples placed in custody room cooler 04/25/03 @ 1440 by Bria

DATA REPORTED ON DRY WEIGHT BASIS

Produced by: Howes, Roberta

Requestor: Brian Farrier

Project Leader: BSTRIGGO

Beginning: 04/22/2003 15:15

RESULTS UNITS

**ANALYTE** 

0.47 U %

% Total Organic Carbon

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. | Reporting limit is an estimate. | N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. | K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.

**EPA - REGION IV SESD, ATHENS, GA** 

Production Date: 06/09/2003 15:36

Sample

6698

FY 2003

Project: 03-0558

Facility: Gulf States Creosoting

Program: SF

ld/Station: GS-08-SD / Media: SEDIMENT

**SPECIFIED TESTS** 

Flowood, MS

Produced by: Howes, Roberta Requestor: Brian Farrier Project Leader: BSTRIGGO Beginning: 04/22/2003 15:50

Ending:

samples placed in custody room cooler 04/25/03 @ 1440 by Bria

DATA REPORTED ON DRY WEIGHT BASIS

**ANALYTE RESULTS UNITS** 

0.41 U % % Total Organic Carbon

U-Analyte not detected at or above reporting limit. | J-Identification of analyte is acceptable; reported value is an estimate. | UJ-Analyte not detected at or above reporting limit. Reporting limit is an estimate. N-Presumptive evidence analyte is present; analyte reported as tentative identification. | NJ-Presumptive evidence analyte is present; analyte reported as tentative identification. Reported value is an estimate. K-Identification of analyte is acceptable; reported value may be biased high. Actual value expected to be less than the reported value.

L-Identification of analyte is acceptable; reported value may be biased low. Actual value expected to be greater than reported value.

NA-Not Analyzed. | NAI-Not Analyzed due to Interferences. | A-Analyte analyzed in replicate. Reported value is "average" of replicates.



# UNITED STATES ENVIRONMENTAL PROTECTION AGENCY REGION 4

Science and Ecosystem Support Division Enforcement and Investigations Branch 980 College Station Road Athens, Georgia 30605-2720

September 09, 2007

4SESD-EIB

# **MEMORANDUM**

SUBJECT: Pipe Tracing Investigation at Gulf States Creosoting site,

Flowood, Mississippi SESD Project #07-0710

FROM:

Brian Striggow

Superfund and Air Section

THRU:

Danny France, Chief

Superfund and Air Section

TO:

Jose Negron

**Emergency Response Branch** 

On September 20, 2007, Brian Striggow of the Science and Ecosystem Support Division (SESD) met with Jose Negron and Alyssa Hughes of the EPA Superfund Division along with Richard Ball and Phillip Weathersby of the Mississippi Department of Environmental Quality (MDEQ) to conduct an investigation at the Gulf States Creosoting Site. The limited purpose of the investigation was to ascertain the existence of a buried pipe system and whether said system could be contributing to contamination found in oxbow lakes adjacent to the site.

Recently, the Consteelco Company, which operates on the southern portion of the Gulf States Creosoting site, discovered apparent creosote contamination on the northwestern portion of their property. The area was investigated by EarthCon Consultants and copies of their report have been forwarded to both EPA and MDEQ. Per the EarthCon report, various creosote-related compounds were found on the northwest corner of the Consteelco property. The report also mentions several broken pieces of large diameter concrete pipe found near the contaminated area, postulated to be part of a site storm drain system.

The owner of property immediately across the levee from Consteelco had brought to the attention of Richard Ball of MDEQ the presence of a manhole cover immediately west of the contaminated area at Consteelco. In an earlier reconnaissance, the cover was lifted revealing a brick valve vault with a valve handwheel visible above the water in the vault. The valve appeared to be approximately 6-8 feet below ground surface.

Also of note, creosote odors have consistently been observed in several areas in and near the oxbow lakes (Creosote Slough) west of the Gulf States Creosoting site. The areas of these chronic odors as well as other features mentioned herein are shown in the attached Figure 1, Pipe Traces.

In the earlier reconnaissance, a wooden structure was discovered on the eastern bank of a drainage ditch that runs north-south along the western edge of the Consteelco property. The structure was approximately 20ft north of the Consteelco property and is on a virtual line extending from a storm drain on the Consteelco property through the broken pipe rubble. The actual path of any storm drains in this area remains unknown.

Work in this investigation was begun by excavating in and around the wood structure at the drainage ditch. This structure was originally thought to be a storm drainage outfall, but upon excavation a flat steel cover was found inside, indicating that the structure was a protective structure for the cover. Figure 2, Cross Section of Tile Pipe, shows these structures as well as subsequently revealed features.

The cover was pried up, revealing what appeared to be a cleanout for an east-west running pipe of approximately 8-10 inches diameter. Dry, gray, medium sand approximately half-filled the pipe at the cleanout. No odors were observed when opening the cover and the soil material in the pipe invert had no odor.

The adjacent ditch invert was probed with a tile probe to locate the point where the pipe crossed the ditch. This area was hand-excavated revealing a glazed clay tile pipe with cemented bell-jointed ends of similar diameter to the pipe at the cleanout. A hole was broken into the top of the pipe at this point for access. There was no soil material in the pipe at this point and again no odors were observed.

A steel fish tape was fed into the pipe for approximately 40 feet in the westerly direction. Utiliquest had been retained for pipe locating work on the site and their representative attached a transmitter to the fish tape and marked the location of the fish tape on the ground surface for the 40 feet that the fish tape extended.

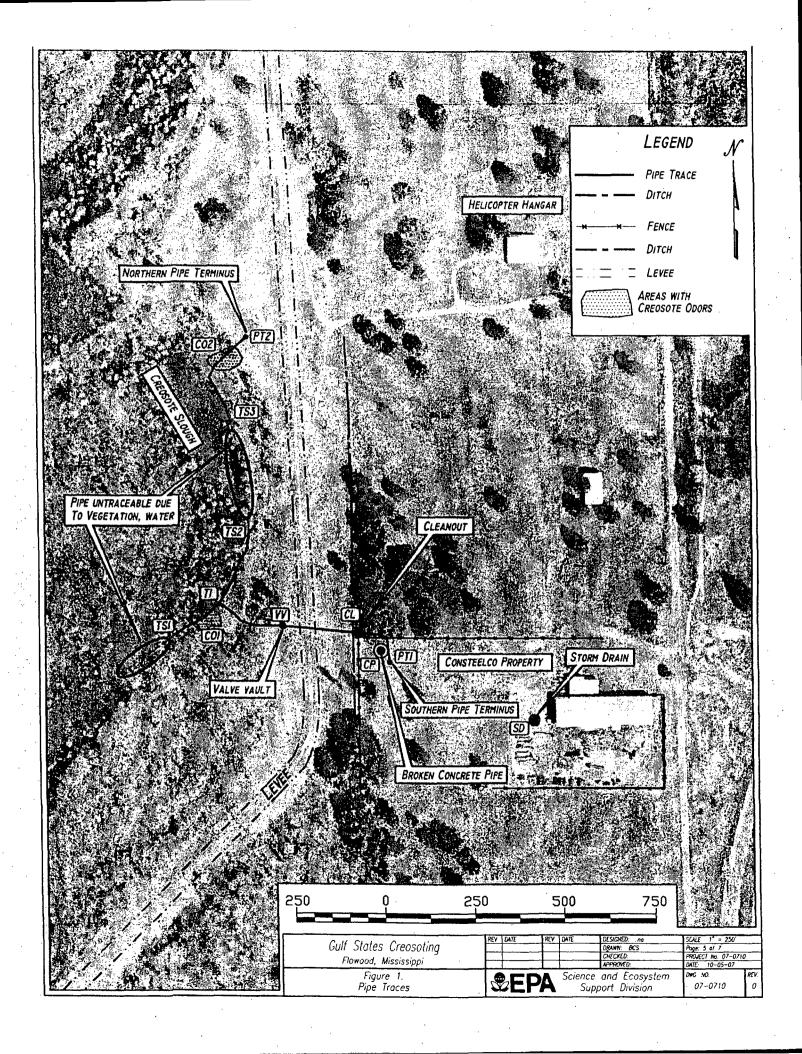
The fish tape was removed and the Utiliquest representative then used inductive-coupled tracing equipment (Pipehorn® brand) to trace the pipe. Inductive-coupled tracing uses a transmitter placed on the ground surface which couples a signal to any nearby underground structures. A handheld receiver is used to detect and trace the signals following the pipes and the ground is marked at the location of pipe trace signals. While clay tile pipe would normally not be expected to conduct a tracing signal, the trace was easily carried to a termination point (labeled PT1 on Figure 1) on the Consteelco property and to the valve vault on the west side of the levee. Of note, the inductive trace directly overlaid the fish tape trace.

The trace was carried west from the valve vault and led to and past an area of creosote odors in creosote slough (labeled CO1 on Figure 1). While the trace appeared to lead further to the west, the trace could not be carried further due to standing water and

The pipe in the bottom of the ditch invert, the cleanout, and the wood structure were destroyed by the excavation. In the event of future work, it should be possible to pick up the trace of the remaining pipe using Pipehorn® or equivalent inductive tracing equipment.

If you have any questions about this document, please contact me at (706) 355-8619 or at email striggow.brian@epamail.epa.gov

Attachment



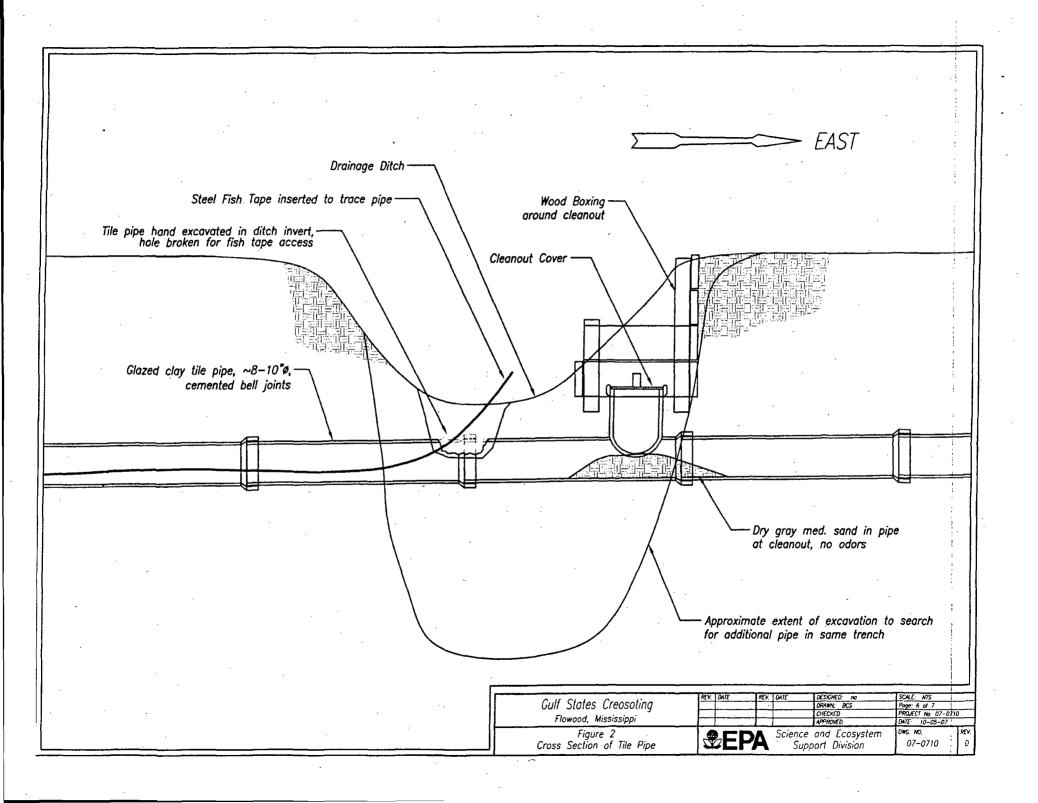


Table 1. Coordinates of Site Features

		WG	584
Map Label	Description	Longitude	Latitude
VV	Valve Vault	-90.14538	32.31057
CO2	Creosote odors, northern	-90.14576	32.31263
CO1	Creosote odors, southern	-90.14608	32.31063
SD	Storm Drain Grating, Consteelco	-90.14315	32.30981
CP	Broken Concrete Pipe at surface	-90.14452	32.31039
CL	Clay Tile Cleanout	-90.14472	32.31050
PT1	Traced Pipe Terminus, southern	-90.14443	32.31028
PT2	Traced Pipe Terminus, northern	-90.14562	32.31277
T1_	Tee west of valve vault	-90.14595	32.31076
TS1	Southernmost stop point of southern pipe trace	-90.14637	32.31052
TS2	Northern stop point of southern pipe trace	-90.14566	32.31128
TS3	Southern stop point of northern pipe trace	-90.14582	32.31223

Reference 19 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423



#### UNITED STATES ENVIRONMENTAL PROTECTION AGENCY

REGION 4 ATLANTA FEDERAL CENTER 61 FORSYTH STREET ATLANTA, GEORGIA 30303-8960

DEC 1 2 2007

Mr. Jerry Banks Mississippi Department of Environmental Quality P. O. Box 10385 Jackson, MS 39289-0385

Subject:

Gulf States Creosoting Company

Flowood, Mississippi

Dear Mr. Banks:

The U.S. Environmental Protection Agency's Emergency Response and Removal Branch (ERRB) conducted a Removal Site Evaluation (RSE) at the above referenced site for potential removal action eligibility under the National Contingency Plan (NCP).

Based on the information collected during the RSE, the On Scene Coordinator (OSC) recommends this site be given a no further action for removal eligibility under EPA's Superfund Removal Program. (See enclosed RSE memo)

This determination does not preclude any other investigation or response action by other parties which may still be appropriate for this site. Should site conditions change or additional information become available, ERRB will re-evaluate this site as necessary.

Should you have any questions concerning ERRB's determination, please contact Jose Negron, OSC at (404) 562-8754, or Jim McGuire, Chief of Removal Operations Section, at (404) 562-8911.

Sincerely,

A. Shane Hitchcock, Chief

Emergency Response & Removal Branch

Enclosure

cc: Dawn Taylor
Tony Moore
Jim McGuire
Jose Negron
Kerri Sanders

# United States Environmental Protection Agency Region IV POLLUTION REPORT

Date: Friday, October 26, 2007

From: Jose Negron, OSC

To: Ralph Howard, EPA Remedial

Subject: Removal Site Evaluation

Gulf States Creosoting

State Highway 468, Flowood, MS

POLREP No.: 1 Site #: A4RF

Reporting Period: D.O. #:

Start Date: 9/20/2007 Response Authority: CERCLA Response Type: Time-Critical

Completion Date: 10/26/2007 NPL Status: Non NPL CERCLIS ID #: MSN000407423 Incident Category: Removal Assessment

RCRIS ID #: Contract #

#### **Site Description**

The former 141-acre Gulf States Creosoting Company is located at 1625 Flowood Drive (Mississippi Hwy 468), Flowood, Rankin County, Mississippi. (The geographic coordinates of the facility are 32'18'43.8" north latitude and 90'58'38.3" west longitude. The facility which operated as a wood treating facility until the mid 1950's is bound by railroad tracks to the north and east, an adjacent business to the south, and marshland/tributary of the Pearl River to the west. The facility is bound by natural barriers and is not fenced. The facility lies within an area comprised of mixed industrial, commercial, and residential uses.

#### **Current Activities**

On September 2007 Brian Striggow from EPA's Science and Ecosytem Support Division along with OSC's Alyssa Hughes and Jose Negron and Richard Ball from the Mississippi Department of Environmental Quality met at the Site to conduct an investigation to ascertain the existence of a buried pipe system that could be contributing to the discharge of creosote sludge onto the oxbow lakes adjacent to the site of the former wood treating facility.

Although an underground pipe network was identified and traced from the site to the marsh there is no evidence of the pipe being a conduit for creosote from the former treatment area onto the marsh. The investigation did not reveal any trace of creosote inside the pipe.

Based on available information it is EPA's determination that there is no clear evidence that establishes that the former wood treating area is the source of creosote at the oxbow. Furthermore ERRB has not identified a migration path that links the suspected source (the former treatment wood treating area) with the presence of creosote at the oxbow.

#### Planned Removal Actions

Based on available information the OSC finds that a time critical removal action at the Gulf States is is not warranted at this time.

**Next Steps** 

ERRB will coordinate with the Site Assesment Branch.

**Key Issues** 

NONE

**Estimated Costs \*** 

	Budgeted	Total To Date	Remaining	% Remaining
Extramural Costs				
Intramural Costs				
			·	
Total Site Costs	\$0.00	\$0.00	\$0.00	0.00%

<sup>\*</sup> The above accounting of expenditures is an estimate based on figures known to the OSC at the time this report was written. The OSC does not necessarily receive specific figures on final payments made to any contractor(s). Other financial data which the OSC must rely upon may not be entirely up-to-date. The cost accounting provided in this report does not necessarily represent an exact monetary figure which the government may include in any claim for cost recovery.

www.epaosc.net/GulfStatesCreosotingFlowoodMS

RSE COORD

12/11/07

City of Flowood - bleinded system Reference 20 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi MSN000407423 - City of Suchson:

166,000 connections

7 surface water

infane

- backup wells for

emergency energy

## **Gulf States**

## Radial Well Search

# **Results Report**

Date of Report: 5/29/2008

#### 0 to 1/4 Mile Buffer

#### **Public Water Supply Wells**

PWS Code PWS System Name	OLWR#	USGS ID# County#	Aquifer Code	Well Status	Elevation	Screentop	Screenbase	Service Connections	System Type
610075-01 CITY OF FLOWOOD-N. WELL	0	121	_	Inactive (1)	. 0	0	0	1500	С

Number of Public Water Supply Wells within 0 to 1/4 mile buffer: 1

1/4 to 1/2 Mile Buffer

#### **Public Water Supply Wells**

PWS Code PWS System Name	OLWR#	USGS ID# County#	Aquifer Code	Well Status	Elevation	Screentop	Screenbase	Service Connections	System Type
610075-02 CITY OF FLOWOOD-N. WELL	0	121	•	Inactive ( I )	0	. 0	0	1500	С

Number of Public Water Supply Wells within 1/4 to 1/2 mile buffer: 1

#### **Source Water Protection Areas**

PWS System ID PWS System Name

610075 CITY

CITY OF FLOWOOD-N. WELL

Number of Source Water Protection Areas within 1/4 to 1/2 mile buffer: 1

1/2 to 1 Mile Buffer

#### Source Water Protection Areas

PWS System ID PWS System Name

610075 CITY OF FLOWOOD-N. WELL

Number of Source Water Protection Areas within 1/2 to 1 mile buffer: 1



## 1 to 2 Mile Buffer

#### **Public Water Supply Wells**

PWS Code PWS System Name	OLWR#	USGS ID#	County#	Aquifer Code	Well Status	Elevation	Screentop	Screenbase	Service Connections	System Type
610034-01 BIERDEMAN-RENTALS	0	K130	121	CCKF	No Data (NS)	275	275	295	110	С
610075-05 CITY OF FLOWOOD-N. WELL	3815	K175	121	SPRT	Active	265	482	562	1500	С
250034-03UNIVERSITY-OF MS MEDICAL CNTR	12890	H194	49	SPRT	=Active	335	680	760	10	Р
250008-01 CITY OF JACKSON	0		49		Surface Water Well	0	0	0	66619	С

Number of Public Water Supply Wells within 1 to 2 mile buffer: 4

#### **Source Water Protection Areas**

PWS System ID	PWS System Name	
610075	CITY OF FLOWOOD-N. WELL	
610017	CITY OF PEARL	•
250034	UNIVERSITY OF MS MEDICAL CNTR	

Number of Source Water Protection Areas within 1 to 2 mile buffer: 3

2 to 3 Mile Buffer

#### **Public Water Supply Wells**

PWS Code	PWS System Name	OLWR#	USGS ID#	County#	Aquifer Code	Well Status	Elevation	Screentop	Screenbase	Service Connections	System Type
610094-01	GULFLINE-INDUSTRIAL-PARK	11319	K076	121	SPRT	No Data ( N3 )	275	706	771	4	. P
610078-01	CAPITAL*56	0		121		Closed	0	0	0	5	N
610075-07	CITY OF FLOWOOD-N. WELL	14933	F070	121	SPRT	Active	275	750	830	1500	С .
<b>1</b> 610017-03	CITY OF PEARL	1300	K120	121	SPRT	Active	315	811	860	7351	С
610017-02	CITY OF PEARL	1302	K119	121	SPRT	Active	320	850	910	7351	<b>C</b> .
√610037-01	CLEVELAND'S TRAILER PARK	0	K093	121	CCKF	Active	267	655	695	135	С
<b>610017-05</b>	CITY OF PEARL	1295	K174	121 .	SPRT	Active	270	550	. 610	7351	C
250034-02	UNIVERSITY OF MS MEDICAL CHTR	3368	H63	49	SPRT	Active	326	694	754	10	P
250034-01	UNIVERSITY-OF-MS-MEDICAL CNTR	3369	H064	49	SPRT	Active	332	716	776	10	Р

Number of Public Water Supply Wells within 2 to 3 mile buffer: 9



#### **Source Water Protection Areas**

PWS System ID	PWS System Name
610075	CITY OF FLOWOOD-N. WELL
610037	CLEVELAND'S TRAILER PARK
610023	CITY OF RICHLAND
610017	CITY OF PEARL
250034	UNIVERSITY OF MS MEDICAL CNTR

Number of Source Water Protection Areas within 2 to 3 mile buffer: 5

Burate well

3 to 4 Mile Buffer

#### <sup>l</sup>Public Water Supply Wells

PWS Code	PWS System Name	OLWR#	USGS ID#	County#	Aquifer Code	Well Status	Elevation	Screentop	Screenbase	Service Connections	System Type
610037-04	CLEVELAND'S TRAILER PARK	0	K188	121	CCKF	Active	265	520	555	135	С
610037-03	CLEVELAND'S TRAILER PARK	0		121	CCKF	Active	260	560	580	` 135	С
610037-02	CLEVELAND'S TRAILER PARK	0	K142	121	CCKF	Active	260	560	600	135	· c
610023-03	CITY OF RICHLAND	14116	K169	121	SPRT	Active ·	265	749	829	2157	С
610017-04	CITY OF PEARL	1303	K095	121	SPRT	Active	315	1093	1143	7351	С
250031-01	PRIMOS NORTHGATE RESTAURANT	0		49		Closed	0.	0	0	. 1	N -
250028-03	JACKSON MUNICIPAL AIRPORT	0	<b>K</b> 199	121	CCKF	Active	310	453	533	1	Р
250028-02	JACKSON MUNICIPAL AIRPORT	0	K28	121	CCKF	Active	320	568	618	1	Р
/6 250028-01 -	JACKSON MUNICIPAL AIRPORT	0	F26	121	CCKF	Active	320	554	614	1	Р

Number of Public Water Supply Wells within 3 to 4 mile buffer: 9

#### **Source Water Protection Areas**

PWS System ID	PWS System Name
610037	CLEVELAND'S TRAILER PARK
610023	CITY OF RICHLAND
610017	CITY OF PEARL
250028	JACKSON MUNICIPAL AIRPORT

Number of Source Water Protection Areas within 3 to 4 mile buffer: 4

fire well

## U.S. EPA REGION IV

# SDMS

Unscannable Material Target Sheet

	Site ID: MSN 000 407423
Nature of Material:	
Map:	Computer Disks:
Photos:	CD-ROM:
Blueprints:	Oversized Report:
Slides:	Log Book:
Other (describe):	deus Map
Amount of material:	

#### **Acronyms**

#### Office of Land & Water Codes AB ABONDONED CHICKEN HOUSE CO COMMERCIAL CATHODIC PROTECTION DO DOMESTIC (HOME) WELL DUPLICATE DW DEWATERING DESTROYED EC **EROSION CONTROL** EXEMPT FΑ FLOW AUGMENTATION FC FISH CULTURE FLOOD PROTECTION FL FΡ FIRE PROTECTION GW GRAVEL WASH INDUSTRIAL IMPOUNDMENT IRRIGATION INSTITUTIONAL LIVESTOCK MUNICIPAL ND NEVER DRILLED NO INFORMATION NOR RENEWED

OBSERVATION WELL OIL WELL OTHER

RECREATIONAL
REMEDIATION WELL
RELIEF WELL
STANDBY
SEDIMENT CONTROL
TEST HOLE
WELL HAS TRANSDUCER

UNUSED

UNKNOWN TEMPORARY YMD1

TEMPORARY YMD2

TEMPORARY YMD3

TEMPORARY YMD4

PRIVATE WATER SYSTEM

RURAL WATER ASSOCIATION

OB OI OT

RE

WA

XX1

XX2

XX3

XX4

.: -	Pul	olic Water Supply Codes
yste	m Type	Description
yste. P	m Type	Description  Non-Transient Non-Community
_	m Type	

		and the second of the second		
Primary Use Description Code				
Α		Air conditioning		
В		Bottling		
C		Commerical		
D		Dewater		
ε		Power generation		
G		Fire protection		
н		Domestic		
J		Irrigation		
J		Industrial (cooling)		
K		Mining		
М		Medicinal		
N		Industrial		
Ρ		Public supply		
Q		Aquaculture (in remarks)		
R		Recreation		
S		Stock supply		
т		Institutional		
Ü		Unused		
Υ		Desalination		
z		Other (explain)		

County #/ FIPS Code	County Name	County #/ FIPS Code	County Name
1	Adams	83	
3	Alcom	85	
5	Amite	87	
7	Attala	89	
9 ,	Benton	91	
11	Bolivar	93	
13	Calhoun	95	
15	Carroll	97	
17	Chickasaw	99	Neshoba
19	Choctaw	101	Newton
21	Claiborne	103	Noxubee
23	Clarke	105	Oktibbeha
25	Clay	107	Panola
27	Coahoma	109	Pearl River
29	Copiah	111	Perry
31	Covington	113	Pike
33	Desoto	115	Pontotoc
35	Forrest	117	Prentiss
37	Franklin	119	
39	George	121	
41	Greene	123	
43	Grenada	125	
45	Hancock	127	
47	Harrison	129	
49	Hinds	131	
51	Holmes	133	
53	Humphreys	135	
55	Issaquena	137	
57	Itawamba	139	
59	Jackson	141	
61	Jasper	143	
63	Jefferson	145	
65	Jefferson Davis	147	
67	Jones	149	
69	Kemper	151	
71	Lafayette	153	
73	Lamar	155	
75	Lauderdale	157	
77	Lawrence	159	
79	Leake	161	. Yalobusha

		Aquifer Codes
quifer Co	ode	Aquiter Name
ALVM:		Alluvial aquiler
CCKF		Cockfield aquifer
COFF		Coffee Sand aquifer
COKR		Coker aquifer
CRNL		Citronelle aquiter
CTHL		Catahoula aquifer
ETMS		Eutaw-McShan aquifer
FRHL		Forest Hill aquifer
GORD		Gordo aquifer
GRMF		Graham Ferry aquifer
HBRG		Hattiesburg aquifer
MSSV		Massive Sand aquifer (lower COKR)
MUWX		Meridian-Upper Wilcox aquifer
MOCN		Miocene aquifer system (includes CTHL, GRMF, HBRG, or PCGL
MRVA		Mississippi River Alluvial aquifer
OLGC		Oligocene aquifer (FRHL not included)
PLZC		Paleozoic aquifer
PCGL		Pascagoula aquiter
RPLY		Ripley aquifer
SPRT		Sparta aquifer system ·
TRCS		
MLCXL		Lower Wilcox aquifer
WLCSN		Middle Wilcox aquifer
WNTL		Winona-Tallahatta aquifer

#### **Disclaimer**

This report produced by the Mississippi Department of Environmental Quality (MDEQ), Groundwater Assessment and Remediation Division. The sources for the data shown are from the Mississippi Automated Resource Information System (MARIS), United States Geological Survey (USGS), and MDEQ. The Mississippi Department of Environmental Quality makes no warranties, expressed or implied, as to the accuracy, completeness, currentness, reliability, or suitability for any particular purpose, of the data contained in this report. In some cases data will be duplicated as the different agencies track and name their sites differently.



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LENGTH		
inches	2.540	centimeters
leet	30.480	centimeters
vard5	- 0.914	meters
mies	1.609	kilometers
9-1909/03/03	44 P.T.	-acre.1

**Reinforcing Steel** 

Wire Mesh

Accessories

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Randy S. Kenner

P.O. Box 6175 1625 Flowood Drive Jackson, MS 39288 Phone: 601-939-3136 Fax: 601-939-0734 E-mail: rsk@consteelco.com

| FGPERALGEL | 10 × 0 F | 30 ) × 585 | 15 × (10 × 1.8) + 32



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DCN: BFW-GSC-0002



Name Gulf States Creosoting Company
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Phone
Project BSte Beconnaissance for Preliminary Assessment / Site Inspection
Preliminary Assessment / Site Inspection

"Rite in the Rain" - a unique all-weather witing surface created to shed water and to enhance the writing mage. Makes it possible to write sharp, legible field data in any kind of weather.

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CONTENTS				
PAGE	REFERENCE	DATE		
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<del>-</del>				

6/3/02
1730 ST EPA Michael Arnett, Brian
Striggow, Don Chris Decker : Shaniela
Pennamon, WESTON-START arrive
@ the former Gulf States Creasating
Company property. The property is
Company property. The property is
& Con Steel Company, Inc.
- Arnett says the creosoting activities —
occurred in the southeastern portion
of the ownerte.
- The owner of the horse farm, -
Jim Webb, said the horse farm
portion of the property underwent-
a Phase I Site Assessment by
Trest Mark Bank
- The horse form has been on the -
property a Tyrs.
- Weldo said the Phase I did not -
raveal anything in the soil or water
- The plant itself comprised = 8 acres
Cof which is now Con Steel Company,
Inc.). The horse form owners own
the property worth of Con Steel Co. Inc.
; south of Con Stoel Co. Inc.

Manager to the state of

The second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second second secon

- The property is > 100 acres in size.

There are goswells on-site which are 1100 to 1200 oldeap. There is a fenced (chain-link) surrounding an observation well for Coke Pipeline (they used to be the owners of the pipeline.

The old revilrand opens that roughlong the horse form were used by rail and which were loaded with cross ties.

- All railroad bed ties are now gone. The horse farm property houses two
helicopters for Channel 3 News. —

- The level that runs along the — costern edge of the property is an - easement

- The property (horse farm: Steel Co.)
is zoned I2 - heavy use industrial.

- The Metal bourn located on the horse fourn porton of the property is used a tractor bourn, it used to be an old storage building during Gulf States Oreosoting Operations.

- The Gulf States Creosoting Operation closed in 1953 or 1954 (according to)

Creasoting bton building: Slough operations were housed in the area where the current Can Steel Co. Inc. is located.

- The old foundation for the former-Gulf States Creosoting building was recently removed. The old foundation was worth of the newbldg: flag pole.

- Con Steel Company is located@

- Love Petrobum owned the Crease Sbugh. Love Petroleum was bought out by a company out of Midland, TX.

- Webb stated people fish in the — Tearl River

-Currently, the majority of the property is covered in green grass.
1900 EPA : START off-orte.

-Note: People in the area obtain their water from the City of Floward.



6/4/02 6/4/02 0830 EPA Miko Amett, Brian Striggau, Chris Decker: START Skameka Tennanan - fish in worthern portion of Creosote Slaigh Lake arrive on site @ Gulf States Steeling - Cut-through Creosole lalo Creasiting: neet w/ state representa-tives Philip Weathersby: Richard Bald, MDFO MDEQ personnel state wells on the property are used for gas storage. The -City of Jackson's surface water intake is just upstracin of the farmer site on the west side of the river (Rear L). — Do recon : go down to Cressoto Stough. Can smell creosote. Clannel which is an entry (maybe) to Cressot Slough -Lake Slough Lake flows w/ Poarl -River. May howe groundwater to -Surface water pathway. - Box turtle on-site Note: Richard Bald States people fish on the other side (west) side of the levee. - Possible eagle wested. (Weathersby) - Commercial fishing on river (cathish)
possibly. (weathersby)

Photo# Description Thoto Mit. Dir 64/02 27 Slagh Lallo βS S Slough Lake 26 Sw W. Portion of Slough balls SP 25 O 24 Overview of Froperty Sp MA N MA SÆ 22 MA NW 21 MA Dranage Ditch MA S Gas well MIA Bock of Consteel Co. SP MA SE Overview of horse faim SP MA Overview of Consteel Co.SP SW 14 Constella barren soil SP NU 14the 13 Upper Channels of Rear River SP

6/4/02 Note: In drainage ditch that rus along the western portion of the property contains "chunks" of Naterial that has petroleum oder. - WLBT Channel & helicopter stored on horsefarm property . -- Several barren aveas on Constello. portion of property-1135 EPA, State; & START person el off-site to go to background. The background will be Jackson-Hep School. rocated on Hwy 25. -1200 Leave Prep School & break for lunch. 1245 Leave lunch area : head to -Sonford Products for recon.



Daniel 1) e/4/02



BILL McGEE Headmaster

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Voice Mail Number: 1-888-527-6891
E-mail: bmcgee@jacksonprep.net



LUKE NEALEY

Director of Finance & Operations

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Telephone: 601-939-8611 Facsimile: 601-936-4068
Voice Mail Number: 1-888-675-9367
E-mail: Inealey@jacksonprep.net

### Late Note:

- Pennamon remembers during oite reconnaissance that no private wells were noticed.
- -On 81 sup 9/12/03, Fennamon Spollo w/ Brion Striggow, EPA SESD who Confirmed that the marsh area adjacent to (west of) the Gulf States Creosoting property flows into the Pearl River.
- There were no residential structures within 200' of the facility.
- On August April 1, 2003, STARTS
  Joe Baer & Shaniera Pennaruon met
  w/ EPA's Michael Arnett, Brian
  Striagow, & Brian Famer. During
  the meeting, sample strategy was
  discussed & it was decided
  in order to determine if any
  contamination was flowing into the
  Poarl River between the

And Market States

background sample & the furthest proposed downstream sample, a Control sample Needs to be collected from Prairie Branch. Any influence from contaminants in Prairie Branch way show up in the intenal between the background sample : the furthest downstream sample (GS-07-3D).

The added sample (to be collected from Prairie Branch) is GS-08-5D.

Reference 22 Gulf States Creosoting Company Flowood, Rankin Co., Mississippi

#### U.S. Census Bureau

State & County QuickFacts

## Rankin County, Mississippi

People QuickFacts	Rankin County	Mississippi
Population, 2006 estimate	135,830	2,910,540
Population, percent change, April 1, 2000 to July 1, 2006	17.8%	2.3%
Population, 2000	115,327	2,844,658
Persons under 5 years old, percent, 2006	6.9%	7.2%
Persons under 18 years old, percent, 2006	24.4%	26.1%
Persons 65 years old and over, percent, 2006	10.5%	12.4%
Female persons, percent, 2006	50.9%	51.6%
White persons, percent, 2006 (a)	79.4%	60.9%
Black persons, percent, 2006 (a)	18.9%	37.1%
American Indian and Alaska Native persons, percent, 2006 (a)	0.2%	0.5%
Asian persons, percent, 2006 (a)	0.8%	0.8%
Native Hawaiian and Other Pacific Islander, percent, 2006 (a)	Z	
Persons reporting two or more races, percent, 2006	0.6%	0.79
Persons of Hispanic or Latino origin, percent, 2006 (b)	1.8%	1.89
White persons not Hispanic, percent, 2006	77.8%	59.39
Living in same house in 1995 and 2000, pct 5 yrs old & over	51.0%	58.59
Foreign born persons, percent, 2000	1.6%	1.49
Language other than English spoken at home, pct age 5+, 2000	3.6%	3.69
High school graduates, percent of persons age 25+, 2000	81.8%	72.99
High school graduates, percent of persons age 25+, 2000 Bachelor's degree or higher, pct of persons age 25+, 2000	81.8% 23.8%	er er er er er er er er er er er er er e
ကြောင်းလေးသောက လက်ပြောက်သောကလေးသည်။ သောက်သည် သည် သည် သည် သည် သည် သည် သည်။ သည် သည် သည် သည် သည် သည် သည် သည် သည်	Caracteria de la companya de la companya de la companya de la companya de la companya de la companya de la comp	16.99
Bachelor's degree or higher, pct of persons age 25+, 2000	23.8%	16.99 607,57
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000	23.8% 18,789	16.99 607,57 24.
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000	23.8% 18,789 25.5	16.99 607,57 24. 1,241,48
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006	23.8% 18,789 25.5 53,180	16.99 607,57 24. 1,241,48 72.39
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000	23.8% 18,789 25.5 53,180 77.1%	16.99 607,57 24. 1,241,48 72.39 13.39
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000	23.8% 18,789 25.5 53,180 77.1% 12.5%	16.99 607,57 24. 1,241,48 72.39 13.39 \$71,40
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600	16.99 607,57 24 1,241,48 72.39 13.39 \$71,40
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089	16.9° 607,57 24 1,241,48 72.3° 13.3° \$71,40 1,046,43
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62	16.99 607,57 24 1,241,48 72.39 13.39 \$71,40 1,046,43 2.6 \$34,27
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000 Median household income, 2004	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62 \$47,580	16.99 607,57 24. 1,241,48 72.39 13.39 \$71,40 1,046,43 2.6 \$34,27 \$15,88 19.3
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000 Median household income, 2004 Per capita money income, 1999	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62 \$47,580 \$20,412	16.96 607,57 24 1,241,48 72.3 13.3 \$71,40 1,046,43 2.6 \$34,27 \$15,88
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000 Median household income, 2004 Per capita money income, 1999 Persons below poverty, percent, 2004	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62 \$47,580 \$20,412 10.6%	16.99 607,57 24.  1,241,48 72.39 13.39 \$71,40  1,046,43 2.6 \$34,27 \$15,85 19.3
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000 Median household income, 2004 Per capita money income, 1999 Persons below poverty, percent, 2004  Business QuickFacts	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62 \$47,580 \$20,412 10.6%  Rankin County	16.99 607,57 24 1,241,48 72.39 13.39 \$71,40 1,046,43 2.69 \$34,27 \$15,85 19.3
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000 Median household income, 2004 Per capita money income, 1999 Persons below poverty, percent, 2004  Business QuickFacts Private nonfarm establishments, 2005	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62 \$47,580 \$20,412 10.6%  Rankin County 3,092 48,386	16.96 607,57 24 1,241,48 72.3 13.3 \$71,40 1,046,43 2.6 \$34,27 \$15,88 19.3  Mississippi 60,54 926,95
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000 Median household income, 2004 Per capita money income, 1999 Persons below poverty, percent, 2004  Business QuickFacts Private nonfarm establishments, 2005 Private nonfarm employment, 2005 Private nonfarm employment, percent change 2000-2005	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62 \$47,580 \$20,412 10.6%  Rankin County 3,092 48,386 21.5%	16.99 607,57 24 1,241,48 72.39 13.39 \$71,40 1,046,43 2.69 \$34,27 \$15,85 19.3  Mississippi 60,549 926,950 -3.19
Bachelor's degree or higher, pct of persons age 25+, 2000 Persons with a disability, age 5+, 2000 Mean travel time to work (minutes), workers age 16+, 2000 Housing units, 2006 Homeownership rate, 2000 Housing units in multi-unit structures, percent, 2000 Median value of owner-occupied housing units, 2000 Households, 2000 Persons per household, 2000 Median household income, 2004 Per capita money income, 1999 Persons below poverty, percent, 2004  Business QuickFacts  Private nonfarm establishments, 2005  Private nonfarm employment, 2005	23.8% 18,789 25.5 53,180 77.1% 12.5% \$98,600 42,089 2.62 \$47,580 \$20,412 10.6%  Rankin County 3,092 48,386	

Geography QuickFacts R	ankin County	Mississippi
Federal spending, 2004 (\$1000)	531,893	22,337,697 <sup>1</sup>
Building permits, 2006	1,562	16,618
Accommodation and foodservices sales, 2002 (\$1000)	135,948	5,486,105
Retail sales per capita, 2002	\$11,448	\$8,724
Retail sales, 2002 (\$1000)	1,392,855	25,017,531
Wholesale trade sales, 2002 (\$1000)	1,303,711	19,215,751
Manufacturers shipments, 2002 (\$1000)	1,401,783	38,276,054
Women-owned firms, percent, 2002	23.1%	25.1%
Hispanic-owned firms, percent, 2002	F	0.7%
Native Hawaiian and Other Pacific Islander owned firms, percent, 2002	F	0.1%
Asian-owned firms, percent, 2002	F	1.6%
American Indian and Alaska Native owned firms, percent, 2002	<b>. F</b> .	0.4%

Geography QuickFacts	Rankin County	Mississippi
Land area, 2000 (square miles)	774.52	46,906.96
Persons per square mile, 2000	148.8	60.6
FIPS Code	121	28
Metropolitan or Micropolitan Statistical Area	Jackson, MS Metro Area	

<sup>1:</sup> Includes data not distributed by county.

Source U.S. Census Bureau: State and County QuickFacts. Data derived from Population Estimates, Census of Population and Housing, Small Area Income and Poverty Estimates, State and County Housing Unit Estimates, County Business Patterns, Nonemployer Statistics, Economic Census, Survey of Business Owners, Building Permits, Consolidated Federal Funds Report

Last Revised: Wednesday, 02-Jan-2008 15:10:38 EST

<sup>(</sup>a) Includes persons reporting only one race.

<sup>(</sup>b) Hispanics may be of any race, so also are included in applicable race categories.

D: Suppressed to avoid disclosure of confidential information

F: Fewer than 100 firms

FN: Footnote on this item for this area in place of data

NA: Not available

S: Suppressed; does not meet publication standards

X: Not applicable

Z: Value greater than zero but less than half unit of measure shown

#### **ENDANGERED SPECIES OF MISSISSIPPI**

MISSISSIPPI NATURAL HERITAGE PROGRAM

- 2002 -

SPECIES NAME	COMMON NAME	GLOBAL RANK	STATE RANK	FEDERAI STATUS
BIVALVIA				
ACTINONAIAS LIGAMENTINA	MUCKET	G5	S1	
CYCLONAIAS TUBERCULATA	PURPLE WARTYBACK	G5	<b>S1</b>	
ELLIPTIO ARCTATA	DELICATE SPIKE	G3G4	<b>S1</b>	•
ELLIPTIO DILATATA	SPIKE -	G5	S1	
EPIOBLASMA BREVIDENS	CUMBERLANDIAN COMBSHELL	G1	<b>S1</b>	(LE,XN)
EPIOBLASMA PENITA	SOUTHERN COMBSHELL	G1	<b>S1</b>	LE
EPIOBLASMA TRIQUETRA	SNUFFBOX	G3	<b>S</b> 1	
LAMPSILIS PEROVALIS	ORANGE-NACRE MUCKET	G2	S1	LŤ
LEXINGTONIA DOLABELLOIDES	SLABSIDE PEARLYMUSSEL	G2	<b>S1</b>	c
MEDIONIDUS ACUTISSIMUS	ALABAMA MOCCASINSHELL	G1 ·	<b>S</b> 1	LT
PLETHOBASUS CYPHYUS	SHEEPNOSE	G3	<b>S1</b>	
PLEUROBEMA CURTUM	BLACK CLUBSHELL	G1	SH	LE ·
PLEUROBEMA DECISUM	SOUTHERN CLUBSHELL	G1G2	S152	LE
PLEUROBEMA MARSHALLI	FLAT PIGTOE	GH	SH	LE
PLEUROBEMA PEROVATUM	OVATE CLUBSHELL	G1	<b>S</b> 1	LE
PLEUROBEMA RUBRUM	PYRAMID PIGTOE	G2	<b>S1</b>	
PLEUROBEMA TAITIANUM	HEAVY PIGTOE .	G1	SH	LE
POTAMILUS CAPAX	FAT POCKETBOOK	G1	·\$1	LE
POTAMILUS INFLATUS	INFLATED HEELSPLITTER	G1	SH	LT
PTYCHOBRANCHUS FASCIOLARIS	KIDNEYSHELL .	G4G5	\$1	
QUADRULA CYLINDRICA CYLINDRICA	RABBITSFOOT	G3T3	<b>S1</b>	
QUADRULA METANEVRA	MONKEYFACE	G4	SH	
QUADRULA STAPES	STIRRUPSHELL	GH	2н	旺
MALACOSTRACA				
FALLICAMBARUS GORDONI	CAMP SHELBY BURROWING CRAWFISH	G1	\$1	. с
INSECTA				o .
NICROPHORUS AMERICANUS	AMERICAN BURYING BEETLE	G2G3	SX	LE
OSTEICHTHYES	•			
ACIPENSER OXYRINCHUS DESOTOI	GULF STURGEON	G312	<b>S1</b>	· LT
CRYSTALLARIA ASPRELLA	CRYSTAL DARTER	G3	<b>S1</b>	
ETHEOSTOMA BLENNIOIDES	GREENSIDE DARTER	G5	SH	
ETHEOSTOMA RUBRUM	BAYOU DARTER	G1	S1	LT
NOTROPIS BOOPS	BIGEYE SHINER	G5	<b>S1</b>	
NOTROPIS CHALYBAEUS	IRONCOLOR SHINER	G4	. S2	
NOTURUS EXILIS	SLENDER MADTOM	G5	51	
NOTURUS MUNITUS	FRECKLEBELLY MADTOM	G3	. 52	
NOTURUS STIGMOSUS	NORTHERN MADTOM	G3	<b>S1</b>	
PERCINA AURORA	PEARL DARTER	GI	<b>S</b> 1	. с
PERCINA PHOXOCEPHALA	SLENDERHEAD DARTER	G5	S1	
PHENACOBIUS MIRABILIS	SUCKERMOUTH MINNOW	G5 `	· \$1	
PHOXINUS ERYTHROGASTER	SOUTHERN REDBELLY DACE	· G5	<b>S2</b>	
SCAPHIRHYNCHUS ALBUS	PALLID STURGEON	G1	S1 .	LE
SCAPHIRHYNCHUS SUTTKUSI	ALABAMA STURGEON	. G1	<b>S1</b>	LE

November 16, 2001

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<sup>\*</sup> Note: Listed plants are only protected federally, no state protection is provided.

SPECIES NAME	COMMON NAME	GLOBAL RANK	STATE RANK	FEDERAL STATUS
ANADINDIA	·			
AMPHIBIA				
AMPHIUMA PHOLETER	ONE-TOED AMPHIUMA	G3	<b>S1</b>	
ANEIDES AENEUS	GREEN SALAMANDER	G3G4	S1	
EURYCEA LUCIFUGA	CAVE SALAMANDER	G5	S1	
GYRINOPHILUS PORPHYRITICUS	SPRING SALAMANDER	G5	51	25
RANA SEVOSA	DARK GOPHER FROG	G1	\$1	PE
REPTILIA				
CARETTA CARETTA	LOGGERHEAD; CABEZON	G3	S1B,SZN	٤T
CHELONIA MYDAS	GREEN TURTLE	G3	SZN	(LE,LT)
DERMOCHELYS CORIACEA	LEATHERBACK; TINGLAR	G2	szn ·	LE
DRYMARCHON CORAIS COUPERI	EASTERN INDIGO SNAKE	G4T3	\$1	LT
ERETMOCHELYS IMBRICATA	HAWKSBILL; CAREY	G3 .	SZN	re
FARANCIA ERYTROGRAMMA	, RAINBOW SNAKE	G5	<b>S2</b>	
GOPHERUS POLYPHEMUS	GOPHER TORTOISE	G3	\$2	(PS:LT) .
GRAPTEMYS FLAVIMACULATA	YELLOW-BLOTCHED MAP TURTLE	G2	\$2	LT
GRAPTEMYS NIGRINODA	BLACK-KNOBBED MAP TURTLE	G3	<b>S2</b>	
GRAPTEMYS OCULIFERA	RINGED MAP TURTLE	G2	52	LT
HETERODON SIMUS	SOUTHERN HOGNOSE SNAKE	G2	SH .	
LEPIDOCHELYS KEMPII	KEMP'S OR ATLANTIC RIDLEY	G1	SIN	LE
PITUOPHIS MELANOLEUCUS LODINGI	BLACK PINE SNAKE	G4T3	52	C
PSEUDEMYS POP 1	MISSISSIPPI REDBELLY TURTLE	G?	<b>S1</b>	
AVE\$	•			
CAMPEPHILUS PRINCIPALIS	IVORY-BILLED WOODPECKER	GH	SX	LE
CHARADRIUS ALEXANDRINUS TENUIROSTRIS	SOUTHEASTERN SNOWY PLOVER	G4T3Q	S2B,SZN	
CHARADRIUS MELODUS	PIPING PLOVER	G3	SZN	(LE,LT)
FALCO PEREGRINUS	PEREGRINE FALCON	G4	SZN	(PS:LE)
GRUS CANADENSIS PULLA	MISSISSIPPI SANDHILL CRANE	G5T1	<b>S</b> 1	LE
HALIAEETUS LEUCOCEPHALUS	BALD EAGLE	G4	S1B,52N	(PS:LT,PDL)
MYCTERIA AMERICANA	WOOD STORK	G4	SZN	(PS:LE)
PELECANUS OCCIDENTALIS	BROWN PELICAN	G4	SIN	(PS:LE)
PICOIDES BOREALIS	RED-COCKADED WOODPECKER	G3	. \$1	LE
STERNA ANTILLARUM ATHALASSOS	INTERIOR LEAST TERN	G4T2Q	53?B	(PS:LE)
THRYOMANES BEWICKII	BEWICK'S WREN	G5	S2S3B,SZN	
VERMIVORA BACHMANII	BACHMAN'S WARBLER	GH .	SXB	LE
MAMMALIA	·			
MYOTIS GRISESCENS	GRAY MYOTIS	G3	SAN	LE
MYOTIS SODALIS	INDIANA OR SOCIAL MYOTIS	G2	SAN	LE
PUMA CONCOLOR CORYIT	FLORIDA PANTHER	G5T1	SH	LE ·
TRICHECHUS MANATUS	MANATEE	G2	SZ	LE
URSUS AMERICANUS	BLACK BEAR	GS	S1	(PS)
URSUS AMERICANUS LUTEOLUS	LOUISIANA BLACK BEAR	G5T2	\$1	LT
ISOETOPSIDA				
* ISOETES LOUISIANENSIS	LOUISIANA QUILLWORT	G3	\$2	Œ
DICOTYLEDONEAE	·			
	PRISTIC POTATO AS AN	G2	<b>S1</b>	ŁŢ
* APIOS PRICEANA	PRICE'S POTATO BEAN	G2 G2	51 52	LI LE
* UNDERA MELISSIFOLIA	PONDBERRY	G2 G2	SH	LE LE
* SCHWALBEA AMERICANA	CHAFFSEED	GZ	חנ	LE

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<sup>\*</sup> Note: Listed plants are only protected federally, no state protection is provided.



#### **Gulf States Creosote**

This Map Is For Advisory Purposes Only



Tuesday, 3 June 2008 09:39

Legend

Flood Hazard Zones



Reference 24
Gulf States Creosoting Company
Flowood, Rankin Co., Mississippi
MSN000407423

Reference 25
Gulf States Creosoting Company
Flowood, Rankin Co., Mississippi
MSN000407423

Project Note			
Date: June 5, 2008	Gulf State Creosoting Company Flowood, Rankin County, Mississippi TDD Number: TNA-05-003-0045		
Organization: T N & Associates, Inc., Reg. 4 EPA START Contract			
Name: Nairimer Berríos-Cartagena Signature: For			
Subject: Gulf State Creosoting HRS (H	lazard Ranking System) Wetland Frontage		
The enclosed information contains:			
	d) for the 15 mile radius surrounding the subject and Wildlife National Wetlands Survey Mapper		
determine the location. The identification of t	Vetlands Survey Mapping website, I was able to he HRS target wetlands which are contiguous to lack of data for the area on Geocortex Internet		
	n Points of Entry (PPE) located at the western ownstream approximately one mile from the site.		
I was not able to extrapolate the information topographic map of the Gulf State Creosotin	n from the USFWS maps onto a 1/50,000 scale ng nor measure the wetlands frontage.		
RESPONS (x) None () Phone call (	SE REQUIRED ) Memo ( ) Letter ( ) Report		
cc: (x)File ()Project Manager (	) Principal Investigator ( ) Other (specify)		

