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BASELINE ECOLOGICAL RISK ASSESSMENT
SALMON SITE
LAMAR COUNTY, MISSISSIPPI

Prepared for

DOE Nevada Operations Office
Las Vegas, Nevada

Prepared by

IT CORPORATION
4330 South Valley View Boulevard, Suite 114
Las Vegas, Nevada 89103

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List of Acronyms and Abbreviations

AEC	U.S. Atomic Energy Commission
AWQC	ambient water quality criteria
BERA	Baseline Ecological Risk Assessment
BOD	biochemical oxygen-demand
CERCLA	Comprehensive Environmental Response, Compensation, and Liability Act
cm	centimeter
CRDL	Contract Required Detection Limit
COC	constituent of concern
COPC	constituent of potential concern
CPOM	Coarse Particulate Organic Matter
CPT	cone penetrometer
CSM	Conceptual Site Model
DOE	U.S. Department of Energy
DNA	deoxyribonucleic acid
DRH	Division of Radiological Health
EEW	Environmental Evaluation Work Plan
EPA	U.S. Environmental Protection Agency
EPC	exposure point concentration
EPC_{aq}	aquatic exposure point concentration
EPC_{sed}	sediment exposure point concentration
ER-L	Effects Range - Low
ft	foot
g	gram
g/day	gram per day
GI	gastrointestinal
GPR	ground penetrating radar
IAEA	International Atomic Energy Agency
IDW	investigation-derived waste
in.	inch
ITAS	International Technology Analytical Services
K_d	soil/water partition coefficient
kg	kilogram
kg/d	kilogram per day

List of Acronyms and Abbreviations *(Continued)*

K_{ow}	octanol/water partition coefficient
kt	kiloton
m	meter
mg/ℓ	milligram per liter
MDA	Minimum Detectable Activity
mg/kg	milligram per kilogram
mg/kg/d	milligram per kilogram per day
MSDEQ	Mississippi Department of Environmental Quality
MTL	maximum tolerance level
NPL	National Priorities List
NTS	Nevada Test Site
pCi/g	picocurie per gram
pC/mℓ	picocurie per milliliter
PCB	polychlorinated biphenyls
RI/FS	Remedial Investigation and Feasibility Study
RBP	Rapid Bioassessment Protocols
SA	Source Area
SGZ	surface ground zero
SS	Salmon Site
TIC	Tentatively Identified Compound
UCL	upper confidence limit
μg/kg	microgram per kilogram
μg/ℓ	microgram per liter
μCi	microcurie

Executive Summary

The Salmon Site (SS), formerly the Tatum Dome Test Site, is located in Lamar County, Mississippi, southwest of the city of Hattiesburg and near the communities of Purvis and Baxterville. It was the site of two nuclear and two gas explosion tests conducted between 1964 and 1970 as part of the U.S. Atomic Energy Commission's (now U.S. Department of Energy [DOE]) Vela Uniform Program to improve the United States capability to detect, identify, and locate underground nuclear detonations. The SS is managed by the DOE Nevada Operations Office.

In October 1964, the U.S. Atomic Energy Commission initiated testing activities at the SS with the detonation of a 5.3 kiloton (kt) nuclear device within the Tatum Salt Dome. This first test, designated the Salmon Event, was conducted at a depth of 826 meters (2,710 feet). The test resulted in the formation of a cavity within the salt dome. A second nuclear test, Sterling (0.38 kt), was detonated in the cavity in December 1968. Two methane/oxygen gas detonations were also conducted in the cavity formed by the Salmon Event. Diode Tube and Humid Water, both equivalent to 0.32 kt, were detonated in April 1970.

A consequence of these testing activities is that radionuclides generated during the testing were released into the salt dome, where they are presently contained. During reentry drilling and other site activities, incidental liquid and solid wastes that contained radioactivity were generated. These wastes resulted in some soil, groundwater, and equipment contamination at the SS. During the 1972 decommissioning of the site, most of the contaminated soil and water were collected and disposed either in the cavity left by the tests or in an injection well. The equipment was decontaminated and transported to the Nevada Test Site for disposal. Nonradioactive wastes were disposed in pits at the site; these pits were subsequently backfilled with clean soil and graded.

The DOE is conducting investigations at the SS as part of a Remedial Investigation and Feasibility Study (RI/FS). The RI/FS is the methodology under the Comprehensive Environmental Response, Compensation and Liability Act (CERCLA) for evaluating hazardous waste sites on the National Priorities List (NPL). The SS is not on the NPL, but DOE has voluntarily elected to conduct the evaluation of the SS in accordance with CERCLA.

As a part of the remedial investigation effort, a Baseline Ecological Risk Assessment (BERA) was conducted at the SS. The purpose of this assessment is to gauge ecological and other environmental impacts attributable to past activities at the former test facility. This assessment was designed to be facility-specific so that all stakeholders could participate effectively.

The purpose of the SS BERA was to

- Describe the observed or potential magnitude of adverse ecological effect(s) at the SS and the primary cause(s) of the effect(s)
- Characterize the ecological consequences of the "no further action" remedial alternative.

Abiotic (soil, sediment, and surface water) and biotic (vegetation and mammalian tissues) media were sampled and analyzed. The site was surveyed and the various species utilizing the aquatic and terrestrial habitats recorded. Sensitive species identified on site were the gopher tortoise and Bachman's sparrow.

Data gathered in the field were used in the problem formulation process to select constituents of potential concern (COPC) and identify ecological receptors. The concerns identified for surface waters were aluminum, arsenic, and radiation sources; for sediments, arsenic, barium, cadmium, lead, manganese, zinc, and radiation sources; and for surficial soils, copper, lead, and radiation sources. Potential radiation sources included tritium and radionuclides detected by gamma spectroscopy (alpha, beta, and gamma omitters) and were taken as COPC in terms of total radiation dose received by an ecological receptor. Potential exposure routes and effects of the COPC to the SS receptors were also identified.

Exposure of ecological receptors was characterized by analyzing the transport and fate of COPC through calculations of exposure point concentrations (EPC) for the various media. The 95-percent upper confidence limit of the mean constituent concentrations for all samples, soil/water partition coefficients (K_d), and bioconcentration factors were considered. Biological tissues analyzed included fish, loblolly pine, a crop plant sample, small mammals, a raccoon, an opossum, and several white-tailed deer. Sampled species were selected as ecological receptors, representing producers and various trophic levels. Species not sampled were also selected, including the sensitive species, and all top consumers represented by the red-tailed hawk and the great blue heron.

Effects to these receptors were characterized by comparison of EPC and tissue concentrations to established criteria, benchmarks, and literature-derived toxicity values. Radiation dose modeled for the mouse, fish, and great blue heron did not exceed established protective criteria. Inorganic COPC in the aquatic ecosystems were eliminated when EPC did not exceed established criteria for the protection of aquatic life. Aluminum exceeded the criteria for surface water in the site samples, and in the reference samples. This is believed to be indicative of high natural levels at the SS. The calculated aquatic EPC for barium and manganese did not exceed values estimated to have adverse effects in freshwater systems; however, residue levels in fish tissue samples were found to exceed reference concentrations. No data were available to determine if these tissue levels were toxic. No acute toxicity was observed in toxicity testing, and most communities appeared unimpaired according to Rapid Bioassessment Protocol metrics.

Terrestrial biota tissue results were compared with values reported in the literature, and it was determined that no adverse effects would be expected due to copper or lead. Calculations made to determine the dose of copper and lead to the red-tailed hawk yielded results well below literature-derived No Observable Adverse Effects Levels. Based on residue concentrations in field collected animals, it was concluded that no adverse effects would be expected for sensitive species or game birds at the SS.

Risk was characterized from these exposure and effects data in the context of uncertainty analysis and potential data gaps. Potential data gaps include the lack of toxicity data for barium and manganese and the lack of game bird and reference-site tissue data. Although localized areas of elevated contamination exist at the site, there is enough evidence to conclude that the assessment endpoints were met in being protective of communities and populations at the SS. The no action alternative would be acceptable unless the desire is to be protective of individual organisms. In that case, removal of isolated areas of elevated contamination should be considered.

1.0 Introduction

1.1 Background

The overall objective of the Salmon Site (SS) Project is to conduct an investigation in accordance with the U.S. Environmental Protection Agency (EPA); Comprehensive Environmental Response, Compensation, and Liability Act (CERCLA); U.S. Department of Energy (DOE) Orders; and Mississippi Department of Environmental Quality guidelines. Although the SS is not listed on the National Priorities List, the DOE has voluntarily elected to conduct the evaluation in accordance with CERCLA. This Baseline Ecological Risk Assessment (BERA) was prepared as part of the CERCLA methodology for characterizing the nature and extent of risks posed by uncontrolled hazardous waste sites. The BERA provides the basis to determine whether remedial actions are necessary and, if so, the justification for performing them at the SS.

1.2 Objectives

The primary objective of this BERA is to determine if the constituents present in environmental media at the SS have caused, or could cause, an adverse ecological impact. Specific objectives of this assessment were to

- Review existing site-related ecological data and, if necessary, perform additional field surveys to identify natural resources and terrestrial and aquatic ecological receptors on and near the site.
- Summarize these data into a description of ecological conditions at the site.
- Develop a conceptual model to identify reasonable site-specific exposure pathways and potential terrestrial and aquatic ecological receptors.
- Select constituents of potential concern (COPC) based on results of chemical analysis of abiotic media samples, using physicochemical and ecologically relevant criteria.
- Perform an ecotoxicological assessment (that includes field measurements and laboratory bioassays) to determine the potential for COPC to induce adverse ecological effects in identified terrestrial and aquatic ecological receptors, then identify constituents of concern (COC) and their potential for inducing adverse effects on ecological conditions at the site.

- Prepare a baseline assessment report that documents the analyses and interpretation of results obtained from the field surveys, laboratory investigations, and ecotoxicological analyses. When adverse impacts are suggested but not quantifiable with available data, identify critical data gaps, define additional data requirements, and make recommendations for additional investigations.

These objectives were met through three phases of investigation that were tailored to the specific circumstances and conditions for the SS. Phase I included a general field survey of the flora and fauna associated with the site, placing particular emphasis on identifying sensitive ecosystems, endangered species and their critical habitats, and species possibly consumed by humans or found in human food chains. Sensitive ecosystems include wetlands, floodplains, wildlife breeding areas, and wildlife refuges. This work was described in reports submitted to the DOE Nevada Operations Office (IT, 1992a and 1993). Phase II investigations included collection of biological tissue and site-related media samples, as well as field monitoring and/or observation, bioassay, and analytical investigations. Phase III involved comparison and evaluation of field data, analytical findings, and bioassay data to relevant criteria and other applicable standards, as available, and culminated in the submission of this BERA report.

1.3 Organization

This report is organized in the following manner, which is consistent with the organization suggested in the *Framework for Ecological Risk Assessment* (EPA, 1992):

Section 1.0 Introduction: Outlines objectives and scope for this assessment and provides general information on the facility being investigated.

Section 2.0 Area Description: Gives a broad-based discussion of ecological and ecologically-related resources and features in the region.

Section 3.0 Problem Formulation: Reviews available information in order to provide an understanding of the current extent of potential problems at the site. Information presented includes available environmental sampling data and identified ecological receptors for each trophic level.

Section 4.0 Exposure Characterization: Characterizes contaminant transport and fate phenomena, identifies specific ecological receptors, and quantifies exposure point concentrations from both primary and secondary exposure pathways.

Section 5.0 Ecological Effects Characterization: Discusses quantitative links between contaminant concentrations and effects in receptors. Literature reviews are the primary source of such dose-response information.

Section 6.0 Risk Characterization: Estimates and describes potential risks to ecological receptors, populations, communities, and ecosystems of interest. An uncertainty analysis is also included in this section.

Section 7.0 References: Contains bibliographic information for Sections 1.0 through 6.0.

2.0 Area Description

2.1 Site Location

The SS is located in Lamar County, Mississippi, near the communities of Baxterville and Purvis. The site is comprised of approximately 595 hectares (1,470 acres) of land in Sections 11, 12, 13, and 14 of the Township 2 North, Range 16 West, St. Stephens Meridian. The test site is rectangular with boundaries that run north-south and east-west. Access to within 1.6 kilometers (one mile) of the site is by way of the Columbia-Purvis County road from Purvis. A network of gravel and sand roads provides access to various locations within the site. A general location map of the area is presented in Figure 2-1.

2.1.1 Site Description

The SS was leased by the government and contractor-operated as part of the U.S. Atomic Energy Commission's (AEC) nuclear weapons testing program. The land is currently owned by the DOE, which also holds the exclusive rights to all subsurface activity at the dome. The test site was decommissioned by the DOE in June of 1972. A granite monument (latitude 31° 08′ 32″ north, longitude 89° 34′ 12″ west) marks the location of surface ground zero (SGZ), which is the ground surface point directly above the nuclear test activities. In 1988, the DOE submitted a Preliminary Assessment of the site to the EPA. This assessment concluded that the residual levels of radioactivity in the vicinity of SGZ were not sufficiently high to warrant listing the former facility as a Superfund site.

2.1.2 Site Maps

Aerial photographs taken at different times during the life of the project were used to assist in locating and defining areas of past DOE activity. The photographs were also used to interpret and evaluate site conditions and activities over time.

From the aerial photographs, archived construction maps, and on-site field work, base maps for the SS were prepared (Figure 2-2). For planning the remedial investigation, the sites on the SS were organized into Source Areas (SA) based on the nature of the activities, geographic proximity, and physical setting. Six SAs have been established at the SS (Figures 2-3 through 2-8) each containing from one to nine individual sites (Table 2-1).

2.2 Sources of Site Historical Data

The SS has been extensively investigated as a part of the DOE activities at the site and monitoring by federal and state agencies. An extensive data base of information on the

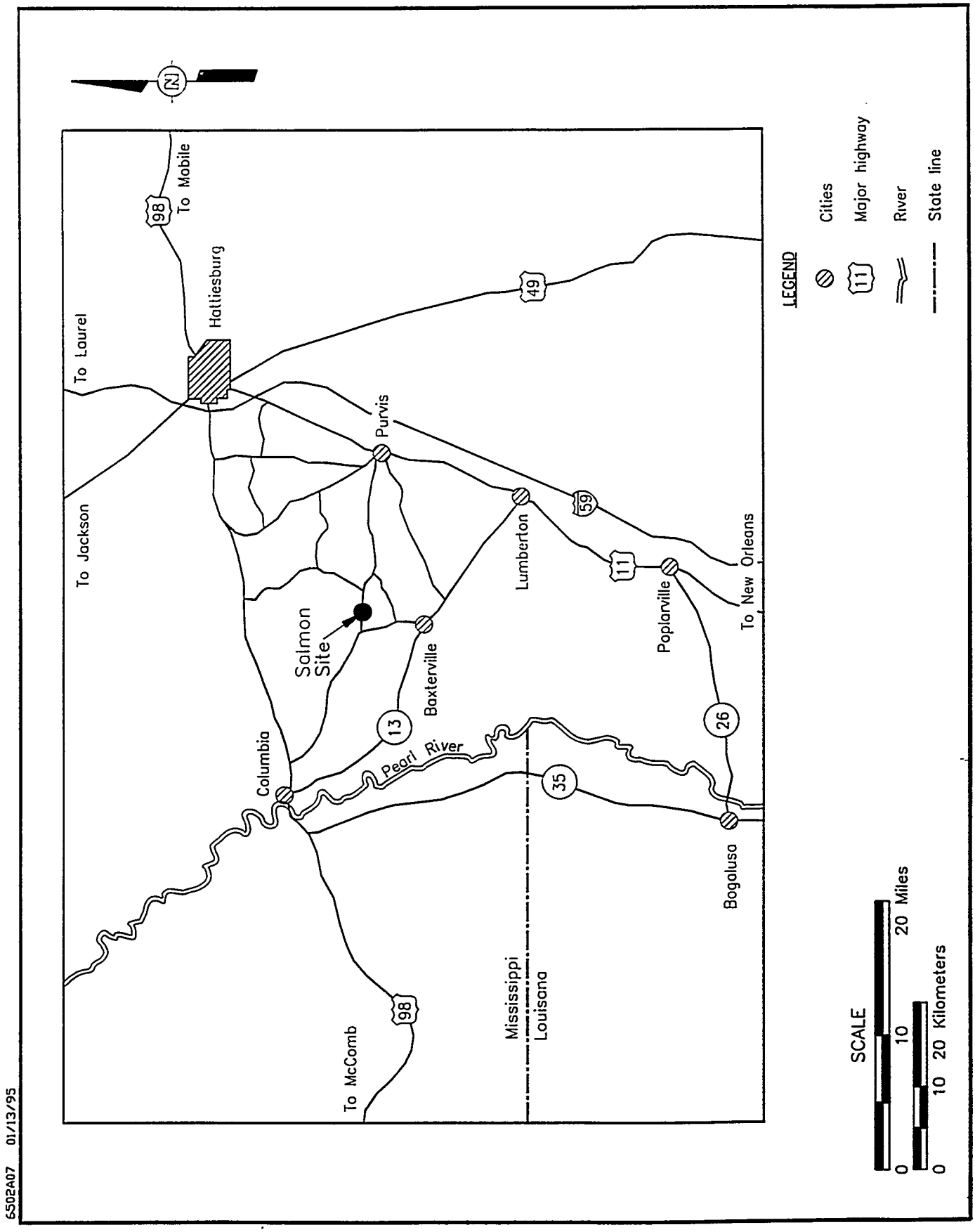

Figure 2-1
General Location Map of the Salmon Site

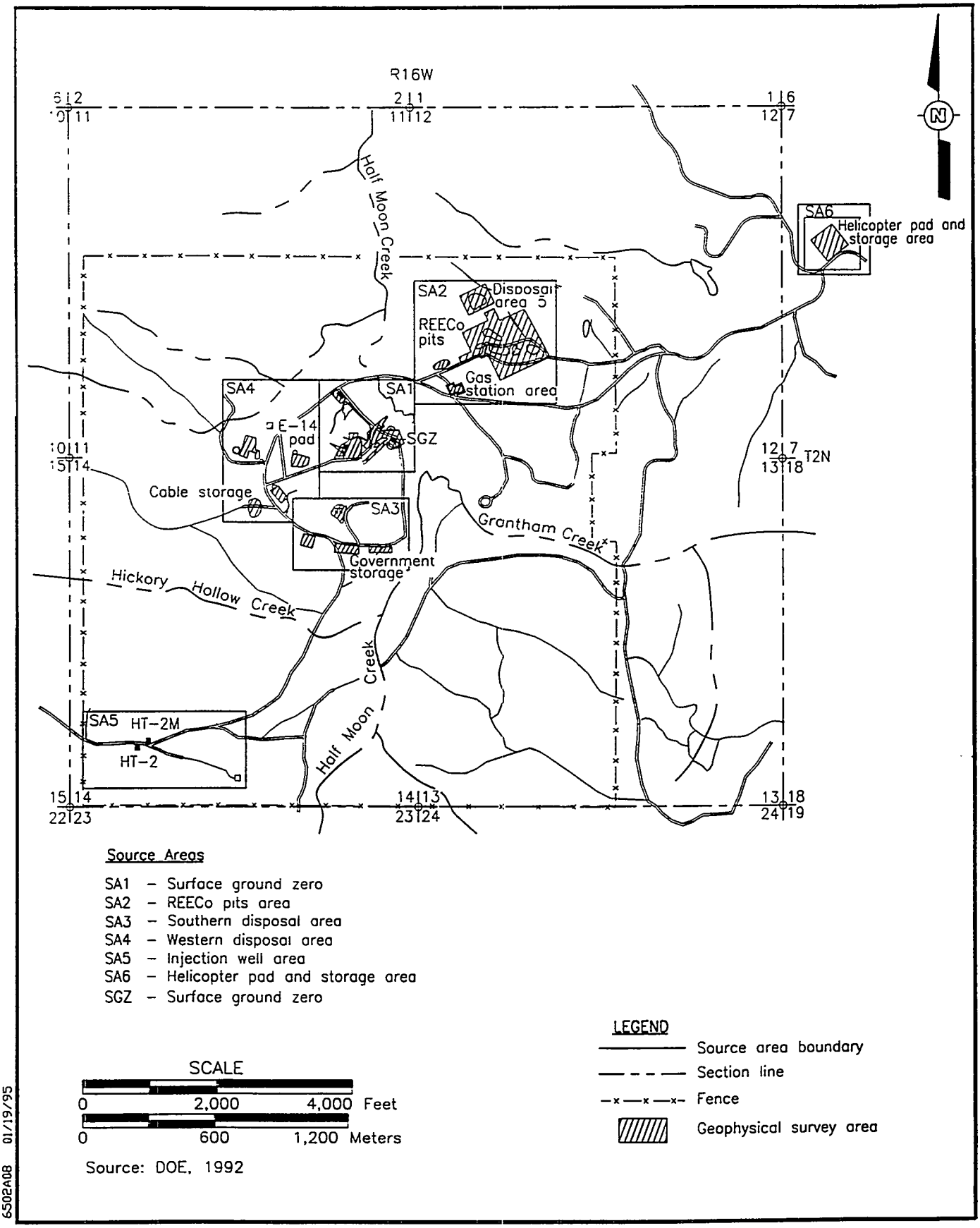

Figure 2-2
Salmon Site Map

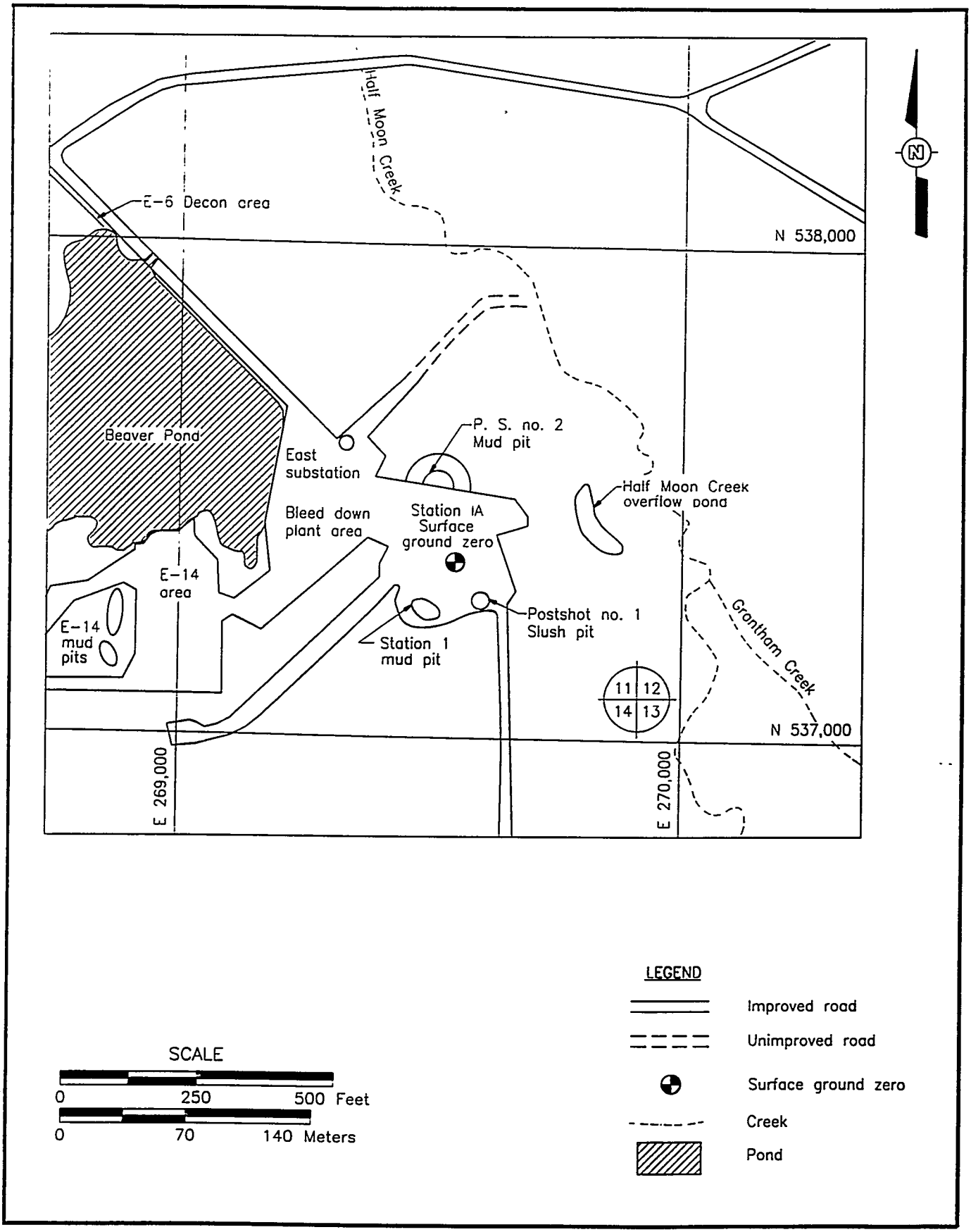

Figure 2-3
Source Area 1
Surface Ground Zero

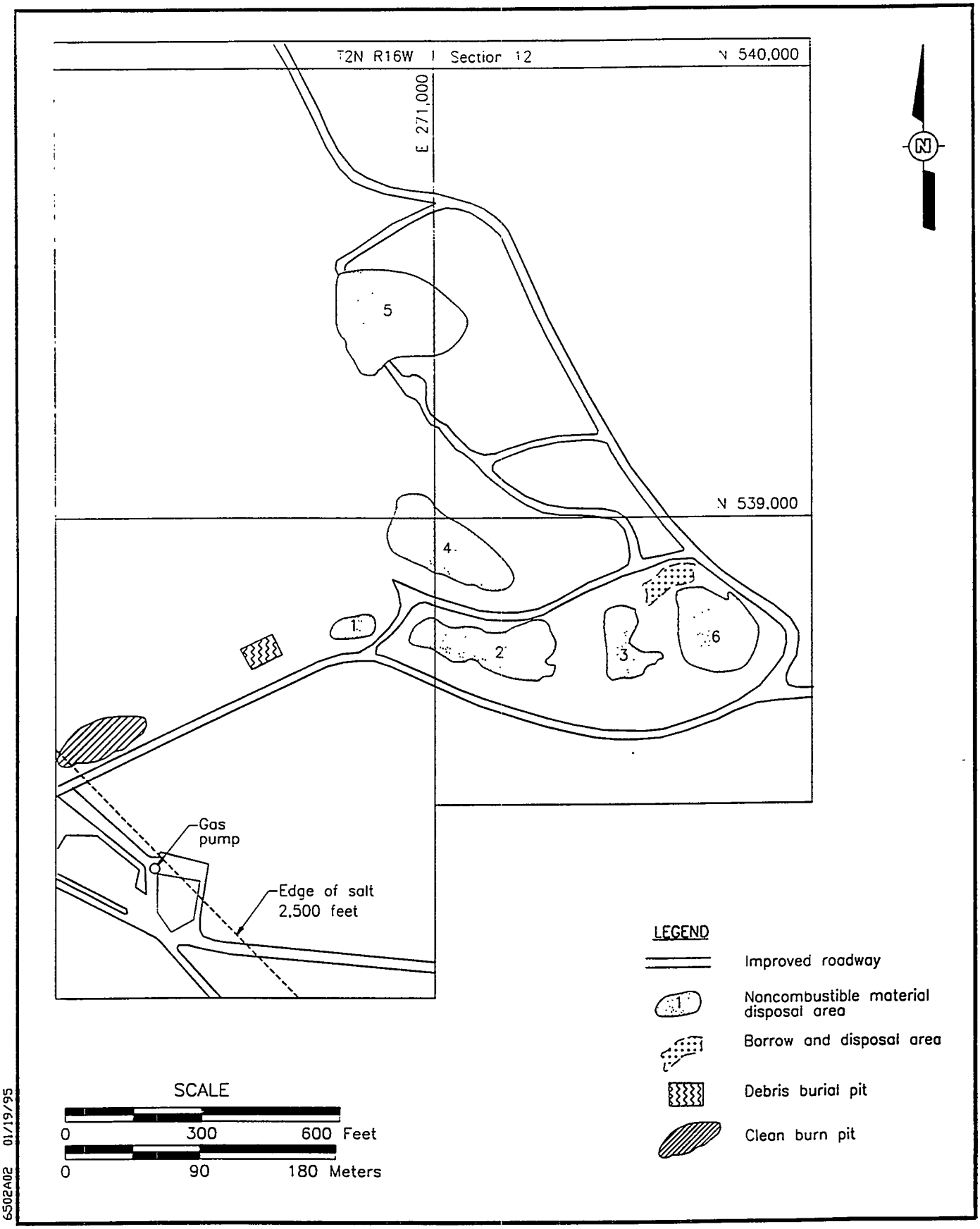

Figure 2-4
Source Area 2
Northern Disposal Area

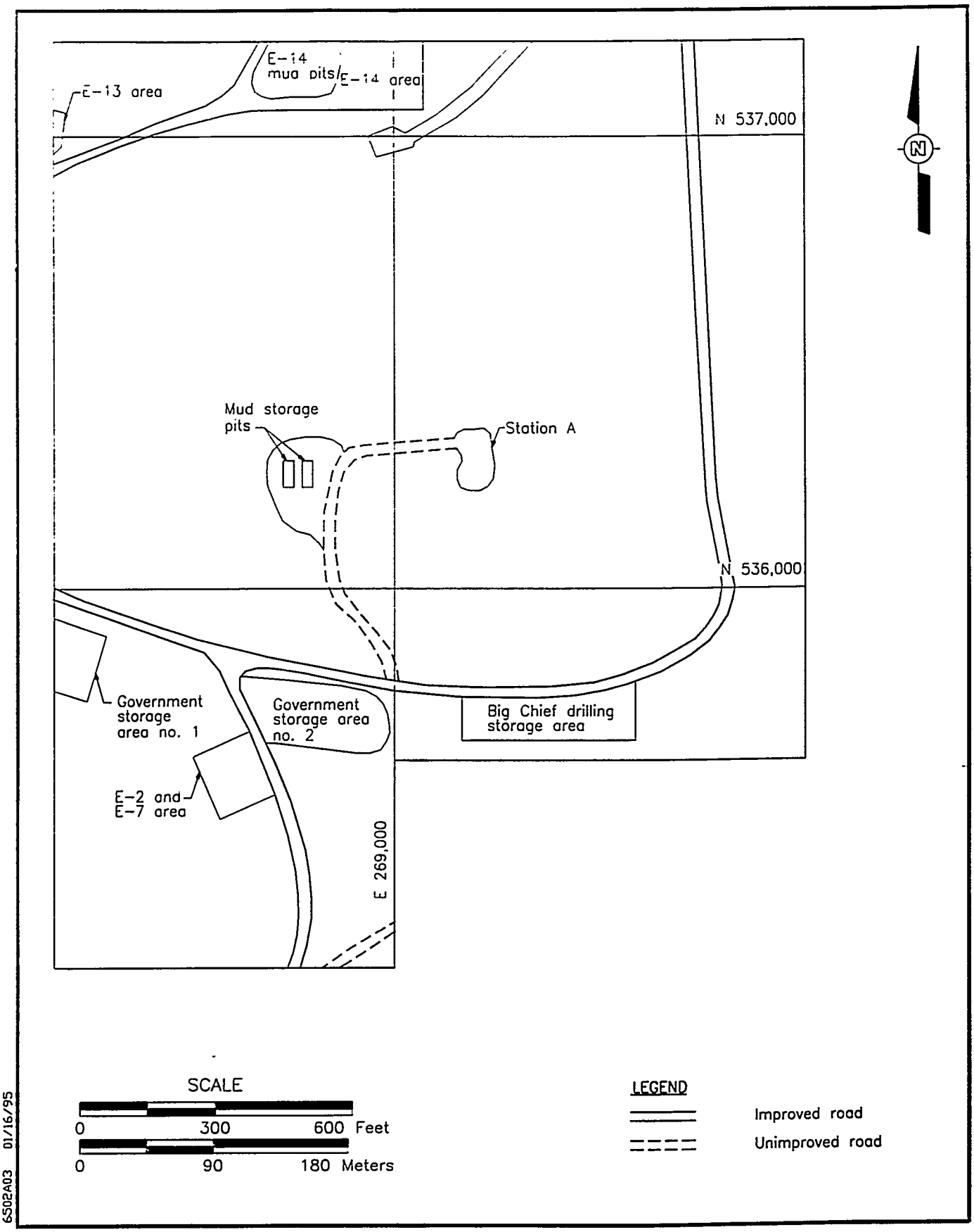

Figure 2-5
Source Area 3
Southern Storage Area

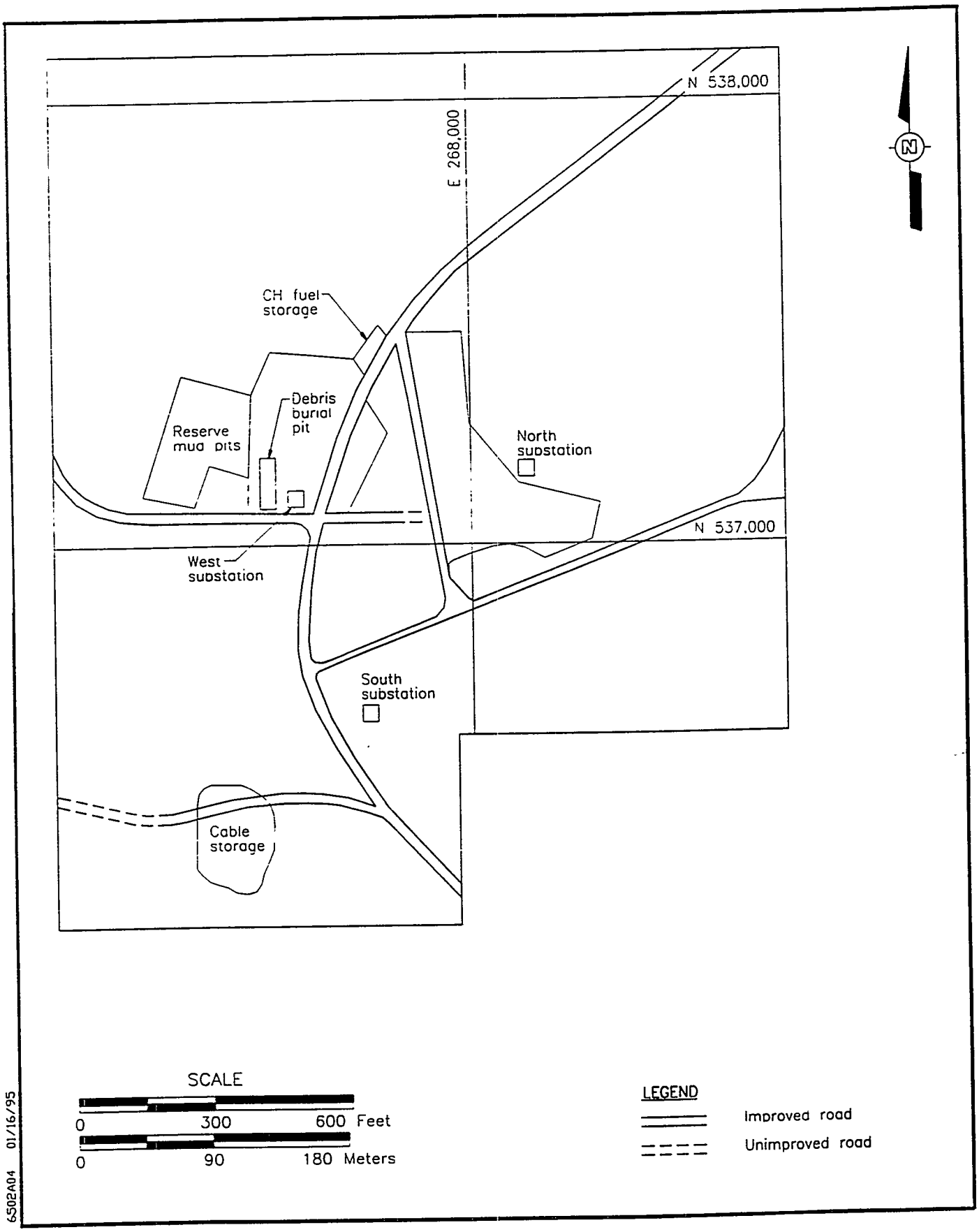

Figure 2-6
Source Area 4
Western Disposal Area

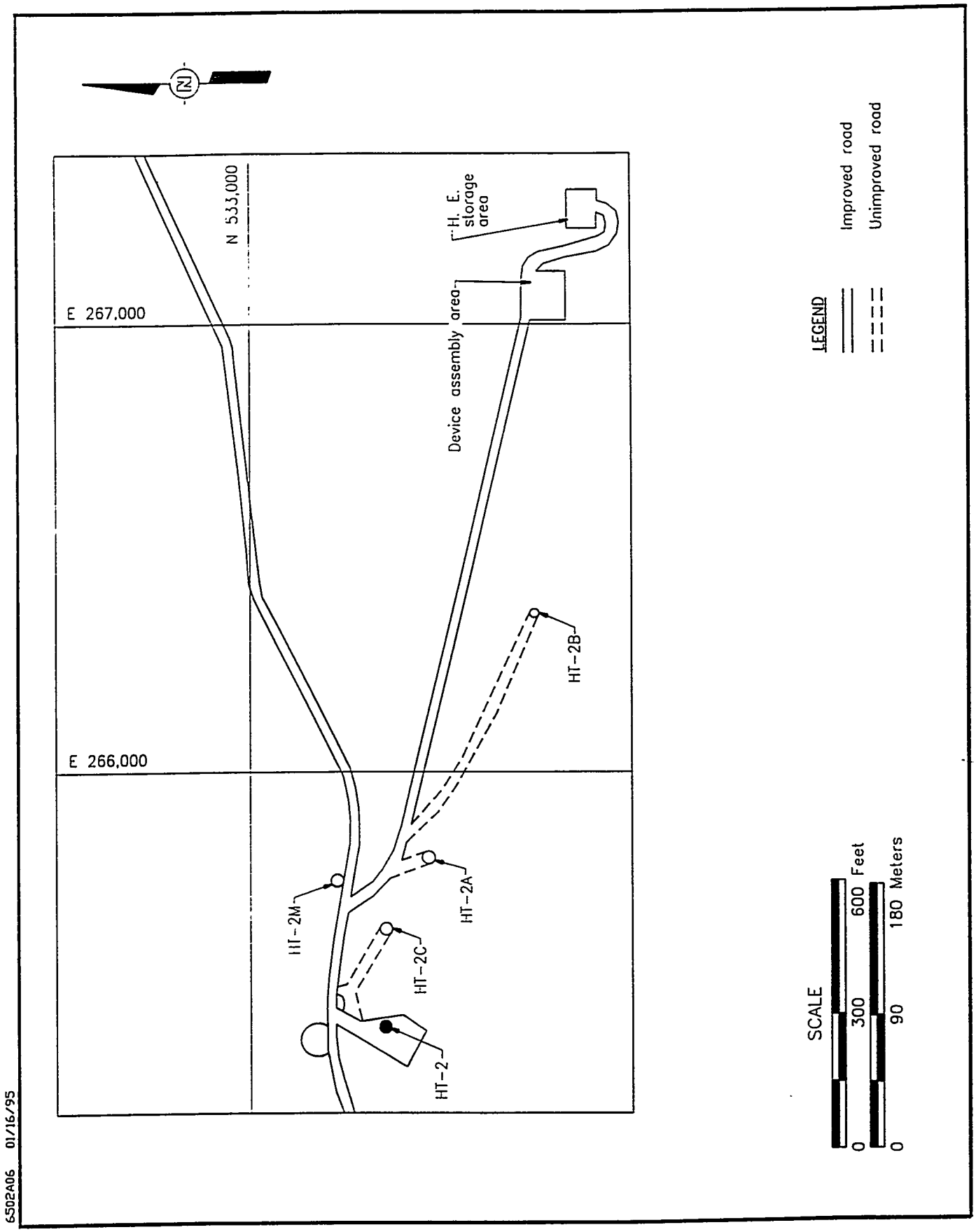

Figure 2-7
Source Area 5
Injection Well Area

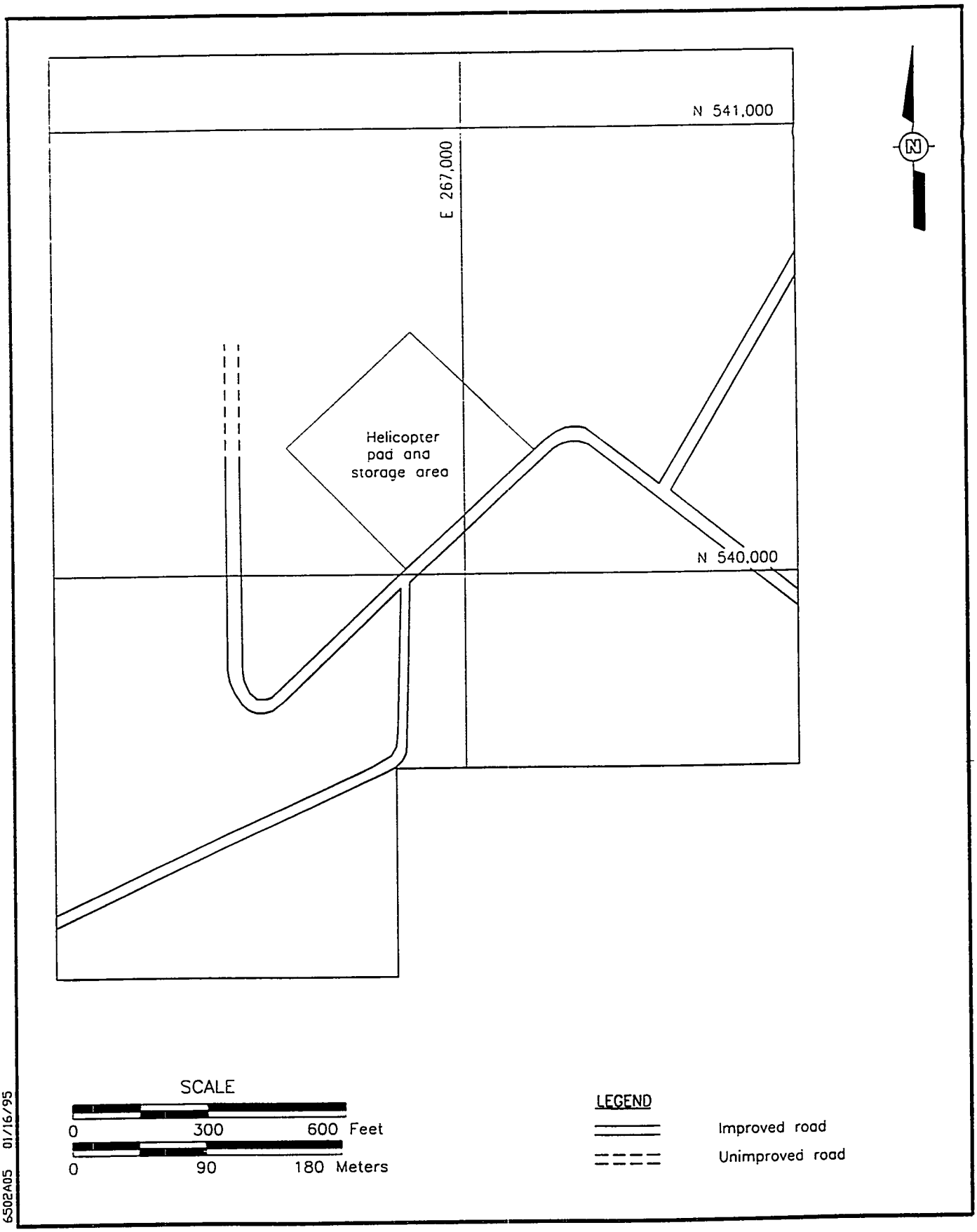

Figure 2-8
Source Area 6
Helicopter Pad and Storage Area

Table 2-1
Sites within each Source Area, Salmon Site

Source Area (SA) 1 - Surface Ground Zero	
1-A	Shot Cavity, Surface Ground Zero
1-B	Beaver Pond
1-C	Half Moon Creek Overflow Pond
1-D	Postshot No. 1 Slush Pit & Mouse Hole
1-E	Bleed-Down Plant Area
1-F	East Substation
1-G	E-14 Pad and Mud Pits
1-H	E-6 Decontamination Pad
1-I	Postshot No. 2 Mud Pit
SA-2 - Northern Disposal Area	
2-A	Reynolds Electrical & Engineering Co., Inc., Disposal Pits
2-B	Clean Burn Pit
2-C	Clean Burn Pit
2-D	Gas Pump
SA-3 - Southern Disposal Area	
3-A	Mud Storage Pits/South Borrow Pit
3-B	Big Chief Drilling Storage Area
3-C	Site E-2 and E-7
3-D	Government Storage Area 1 (Drilling Storage Yard)
3-E	Government Storage Area 2
SA-4 - Western Disposal Area	
4-A	Reserve Mud Pits
4-B	Debris Burial Pit
4-C	West Substation
4-D	CH Fuel Storage
4-E	Cable Storage
4-F	South Substation
4-G	North Substation
SA-5 - Injection Well Area	
5-A	HT-2 Well Area
5-B	HT-2M Well Area
SA-6 - Helicopter Pad and Storage Area	
6-A	Helicopter Pad and Storage Area

operation, waste generation, disposal activities, and physical setting of the site is available. Historical environmental data collected from the early 1970s through 1990 on the SS are chronicled in the following reports:

- *Cleanup Summary Report, Tatum Dome Test Site, Mississippi*, June 1972 (NVO-129) (DOE, 1972)
- *Long-Term Radiological Surveillance Program, Tatum Dome Test Site*, Revised March 1975 (NVO-143) (DOE, 1975)
- *Special Study: Tatum Dome Test Site Lamar County, Mississippi*, Final Report, October 1978 (NVO-200) (DOE, 1978)
- *Animal Investigation Program, Mississippi Results* (EPA, 1974)
- *The Tatum Dome Project, Lamar County, Mississippi* (NVO-225) (DOE, 1980)
- *Onsite and Offsite Environmental Monitoring Report: Radiation Monitoring Around Tatum Salt Dome, Lamar County, Mississippi* (EPA, 1990)
- *Tatum Salt Dome Radiological Monitoring*, Annual Report, March 1990 (MDOH, 1990).

2.2.1 Site Operational History and Current Status

In 1961, the AEC (now the DOE) conducted exploratory drilling, surveying, and road construction at a site located on the Tatum Salt Dome in Lamar County, Mississippi (AEC, 1972). The Tatum Salt Dome was selected as the site of the AEC's Vela Uniform Program designed to determine if a salt dome could be used to mask the seismic signature of underground nuclear tests. Construction for the Salmon test started in early 1963 with the boring of an emplacement shaft into the salt dome. In October 1964, the Salmon device was detonated at a depth of 828 meters (m) (2,716 feet [ft]). The Salmon event had a reported yield of 5.3 kilotons (kt) and resulted in the formation of a cavity with a radius of approximately 33 m (108 ft) within the salt dome (AEC, 1972). Following the event, the point of detonation was investigated by drilling back into the cavity and sampling the contents. To reduce the risk of radioactivity being released from the cavity to the environment during the re-entry drilling, a bleed-down plant was constructed. Gases released during the re-entry boring were routed to the bleed-down plant where they were filtered to reduce the radioactivity to regulatory standards prior to release to the atmosphere.

A second nuclear event, Sterling (0.38 kt), was detonated in December 1966 in the cavity formed by the Salmon Event. Both events were fully contained so there was no uncontrolled

venting of gases or particulates to the surface. There was a minor release of gaseous, radioactive materials during cable cutting and sealing of the holes for the Sterling event (DOE, 1971). It is expected that any such releases would have been quickly dissipated to the atmosphere.

Two methane/oxygen gas detonations were also performed in the cavity left by the Salmon Event. The two tests, Diode Tube and Humid Water, were part of the Miracle Play Series events, both had yields of 0.32 kt, and were detonated in February 1969 and April 1970, respectively.

Following the Humid Water event the DOE began clean-up, restoration, and decommissioning activities at the site.

As a consequence of the testing activities, radionuclides were released into the salt dome where they are presently contained. During re-entry drilling and other site activities, incidental liquid and solid wastes were generated that contained radioactivity. These wastes resulted in some soil, groundwater, and equipment contamination at the site. Most of these wastes and contaminated soil and water were cleaned up and disposed of, either in the shot cavity or in an injection well. The equipment was decontaminated and transported to the Nevada Test Site (NTS) for disposal. Nonradioactive wastes were disposed in pits at the site, which were backfilled with clean soil and graded. The DOE plugged 34 of the 39 exploratory and monitoring wells that were constructed between 1961 and 1972. The site was officially decommissioned on June 30, 1972.

Environmental monitoring at the SS was initiated in 1964 by the Public Health Service and continued until May 1970. Long-term hydrologic monitoring was initiated in 1972 by the EPA and continues to the present. As a part of this monitoring, all potable aquifers, many individual water wells, public water supplies, and some surface waters have been and are currently sampled annually in the vicinity of the SS. The MSDEQ, Division of Radiological Health (DRH), also conducts a quarterly monitoring program.

In addition to the monitoring program maintained by the EPA, the DOE is currently conducting a Remedial Investigation and Feasibility Study (RI/FS) (IT, 1992b) that includes this BERA. Preliminary activities started in July 1992, but the major intrusive tasks could not be conducted until DOE owned the land. The following outlines activities that have been conducted on site since July 1992:

July 1992 Initiation of surveys and RI/FS Work Plan:

Floodplain and Wetland Investigation
- Wetlands delineation

National Environmental Policy Act Surveys
- Avian
- Mammalian
- Reptile/amphibian (gopher tortoise identified on site)
- Vegetation
- Historical/archaeological

Aquatic Surveys
- Rapid Bioassessment Protocols
- Established surface water/sediment sampling stations

Initial Geophysical Investigation
- Established grids and collection of magnetometer data

October 1992

Gopher Tortoise Survey
- Located burrows and habitat evaluation

Continuation of Geophysical Survey
- Cleared survey grids
- Ground penetrating radar (GPR)
- Magnetometer

February 1993

Sample Collection
- Surface water
- Sediment
- Biota
  - Fish, mammals, aquatic invertebrates
- Expanded geophysical survey
  - Line cutting, GPR, magnetometer

April 1993

Soil/Vegetation Sampling
- Sampled SAs for soil and vegetation
- Continued geophysical survey
  - Line cutting, GPR, magnetometer

August 1993

Continuation of RI/FS Work Plan
- Expanded geophysical investigation by 100 percent
  - Clearing lines, GPR, magnetometer
- Clearing/grubbing of areas to be used for investigation-derived waste (IDW) Storage and Decontamination Pad

September 1993
- Construction of new bridge/bypass on Half Moon Creek
- Repaired and widened site roads
- Continued geophysical survey
- Installed fence around decontamination pad

October 1993
- Completed geophysical surveys
- Flagged well and cone penetrometer (CPT) locations

November 1993
- CPT
  - Cleared 18 sites
  - Stratigraphy logs
  - Sampled groundwater and soil with the CPT rig
- Installed 2 wastewater tanks at the decontamination pad

December 1993
- Trench Excavations
  - Dug 17 trenches in Reynolds Electrical & Engineering Co., Inc., pits to locate the magnetic anomalies identified through geophysical surveys
  - Soil sampled from around buried anomalies
- Trench Closure
- Cleared and prepped two drill pads at SGZ

January 1994
- Site Closure
  - Moved waste/drums off site
  - Moved all equipment to office compound
  - Sampled wastewater tank
  - Closed office trailer and demobilized IT Corporation personnel as of January 21

2.2.2 Hazardous Substance and Waste Management

Both radioactive and hazardous wastes were generated as part of the normal operations at the SS. During operation, more than 100 people worked at the SS requiring fuel, electricity, sanitation, waste disposal, and the storage and use of hazardous materials. In addition, following each of the test events, reentry holes were drilled into the shot cavity to determine the subsurface effects of the explosion. These drilling operations generated the largest volume of residual waste materials, primarily contaminated drill cuttings and fluids.

Many of the individual sites within the SAs were investigated during site decommissioning, and contaminated soils and equipment were consolidated and disposed. Radioactively- contaminated equipment was decontaminated and shipped to the NTS for disposal. Radioactively-contaminated soil and drilling fluids were transported to the SGZ area and injected into the shot cavity. Radioactively-contaminated water was also injected into the deep brine aquifer used for

the disposal of oil field brines. Nonradioactively-contaminated wastes were primarily disposed by landfilling them on site.

The initial activities completed under the RI/FS at the SS produced some IDW. These are stored on site pending completion of the land acquisition. Construction debris, such as concrete, steel landing mat, and old culverts from the Half Moon Creek bridge bypass construction was piled to either side of the new bridge. These materials have been surveyed for gross radiological contamination, and samples were analyzed at a laboratory for radiation with negative results. All trees that were removed in the process of building the IDW storage area were piled to the east of the main road near the decontamination pad. These trees are creating a potential biological hazard because they are infested with bark beetles that may attack the live trees along the edges of the IDW storage area. Within the decon-pad compound there are approximately 4,500 gallons of stored waste water from decontamination of the CPT equipment. Laboratory analyses indicate no radiological or hazardous COCs in the water.

2.3 Issues of Public Concern

At the present time, members of the surrounding community are concerned that past activities at the salt dome may pose a risk to the environment and public health. Concerns to the BERA include the following:

- Is radionuclide contamination present in the on-site environmental media?
- Is nonradionuclide contamination present in the on-site environmental media?
- If present, what is the nature and extent of contamination?

Several long-term monitoring programs have been funded by DOE and conducted by EPA and MSDEQ/DRH. In the 18 plus years since these programs were initiated, no radioactivity above naturally occurring background levels has been detected in any of the off-site water supplies. Studies conducted by the MSDEQ/DRH and EPA have found no unusually high incidence of cancer in humans within Lamar County.

3.0 Problem Formulation

Problem formulation is the first step of the ecological risk assessment process. Problem formulation can be defined as a systematic planning step that identifies the major factors to be considered in a particular assessment (EPA, 1992). It establishes the goals, breadth, and focus of the assessment and is linked to the regulatory drivers and policy context of the assessment. The process begins with the initial characterization of the exposure and ecological effects, including evaluating the stressor characteristics, the ecosystem potentially at risk, and the ecological effects expected or observed. Problem formulation describes the relationships among assessment and measurement endpoints, data required, and methodologies that will be used to analyze the data.

3.1 Site Description

This section describes and discusses the physical and biological characteristics of the SS. These characteristics (i.e., geology, water resources, climate, existing use, and biota) provide the building blocks from which the conceptual model of the site was developed.

Overall, the SS has largely returned to a natural state and, except for the monument erected by the DOE at SGZ, there is little indication of the past activities at the site. The areas where soils were excavated have been back filled and seeded and now have a well established cover of vegetation. Numerous species of flora and fauna are abundant at the former test site, and the area is currently used for timber production and hunting.

3.1.1 Physical Features

The SS is situated in the low hills of the pine woods area of the Gulf Coast region. This area is characterized by narrow, flat-topped ridges and intervening valleys that trend predominantly in a south-southeast direction toward the Gulf of Mexico. The maximum relief on site is approximately 30 m (100 ft) with elevations ranging from 75 to 110 m (246 to 360 ft) above mean sea level. Surface ground zero lies east of Half Moon Creek, the principal drainage from the site.

The SS is situated on the northern edge of the Mississippi Salt Basin. The geologic strata deposited in the basin are typically interbedded, discontinuous sandstones and shales with local marls, evaporites, and limestones. Regionally the sediments dip gently to the south and southeast. Locally, the dip may be modified by the intrusion of salt domes.

The SS is located in the Atlantic and Gulf Coastal Plain Groundwater Region. Groundwater occurs throughout the region in a vertical series of aquifers, with the surficial aquifer generally located less than 30 m (100 ft) below surface. On a regional scale, groundwater flow is primarily in a southerly or southeasterly direction. However, the hydrogeologic conditions on the local, near-dome scale, are very complex because of the effect of the dome. At SGZ, the alluvial aquifer extends from near ground surface to a depth of about 12 m (39 ft) below the land surface.

Surface drainage at the SS is through three creeks: Hickory Hollow, Grantham, and Half Moon Creeks. These creeks receive groundwater discharge in the dry season from the alluvial aquifer, which may transport contamination along with it. Tritium and trace levels of organic constituents have been detected in groundwater from the alluvial aquifer near SGZ. Ponds located on the SS also represent important hydrologic elements. The ponds of most concern with regard to contaminated groundwater discharge are the Beaver Pond and Half Moon Creek Overflow Pond (Figure 3-1).

3.1.2 Biological Features

The SS is located within a belt of pine forest that extends along the coastal plains of the South Atlantic and Gulf States. The pine forest is a seral stage of the temperate deciduous forest. These forests can be classified into either one of two major types: loblolly/short leaf and slash/long leaf pine. The continued dominance of the pine species is dependent upon periodic fires and cutting, without which the pines would be replaced by hardwoods, such as oaks (*Quercus* sp.), hickory (*Carya* sp.), and magnolia (*Magnolia* sp.). Most of the SS is open, long-leaf pine forest with a shrub understory.

The diversity of vegetation and perennially available water provide suitable habitat for whitetailed deer (*Odocoileus virginianus*); armadillo (*Dasypus novemcinctus*); eastern cottontail rabbit (*Sylvilagus floridanus*); coyote (*Canis latrans*); raccoon (*Procyon lotor*); opossum (*Didelphis marsupialis*); beaver (*Castor canadensis*); squirrel (*Tamiasciurus hudsonicus* and *Sciurus carolinensis*); several species of small rodents; and numerous bird, reptile, and amphibian species. The small creeks and ponds on site provide limited habitat for large fish species, but do support populations of minnows, which provide forage for egrets, herons, and other piscivorous species. A complete listing of all flora and fauna observed on site is presented in Tables 3-1 through 3-6.

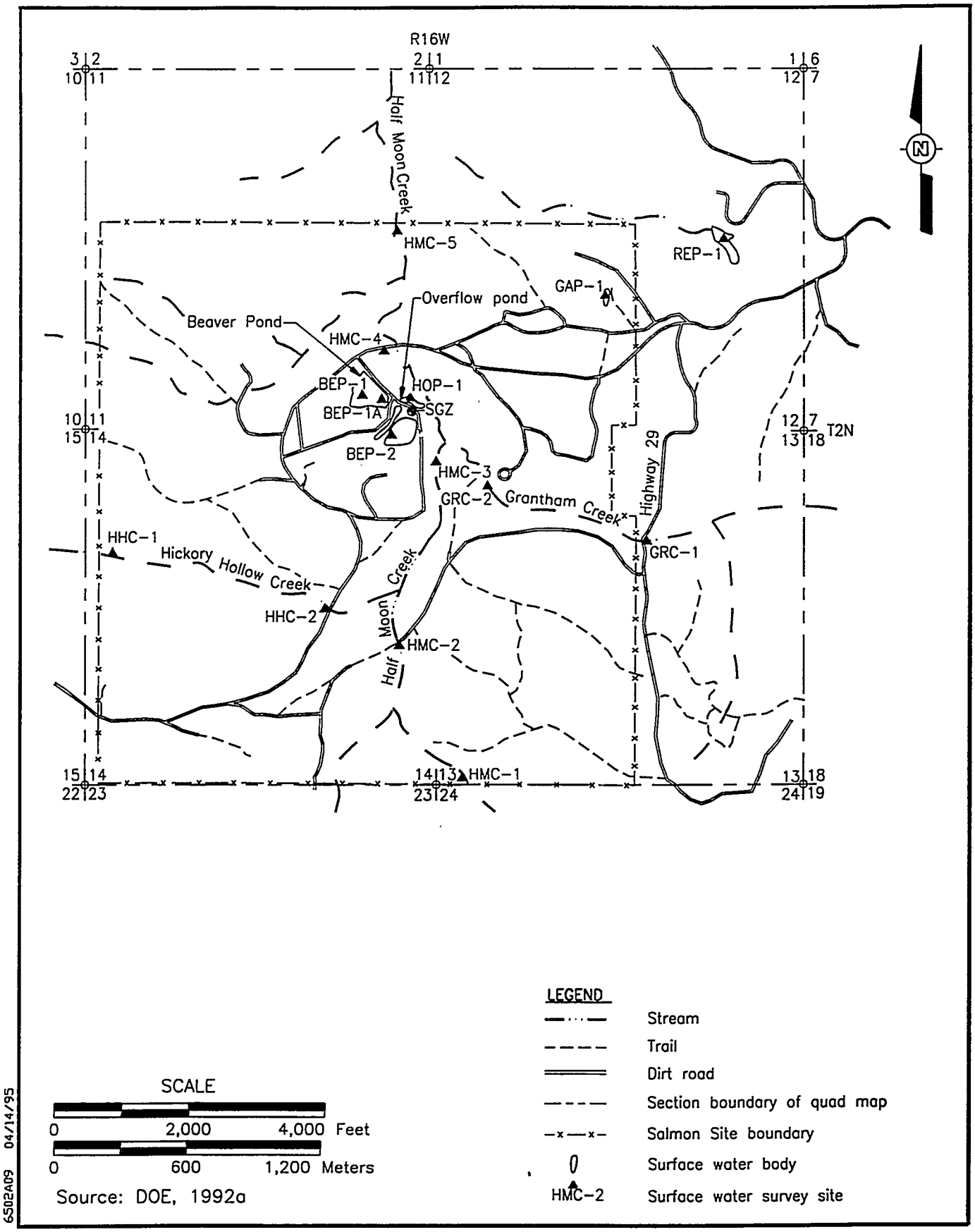

Figure 3-1
Salmon Site Surface-Water and Sediment Sampling Locations

Table 3-1
Species List - Vegetation Survey,
July 13 to 19, 1992,
Salmon Site
(Page 1 of 3)

Common Name	Scientific Name
Sensitive Fern	**Polypodiaceae** *Onoclea sensibilis*
Green Bristlegrass Bristlegrass Foxtail Grasses Cord Grass	**Gramineae** *Setaria viridis* *Setaria* sp. *Alopecurus* sp. *Gramineae* sp. *Spartina pectinata*
Bladder Sedge Umbrella Sedge	**Cyperaceae** *Carex intumescens* *Cyperus strigosus*
Canada Rush Toad Rush	**Juncaceae** *Juncus canadensis* *Juncus bufonius*
Fragrant Water Lily	**Nymphacaceae** *Nymphaea odorata*
Common Greenbriar Long-Stalk Greenbriar Indian Cucumber Root	**Liliaceae** *Smilax rotundifolia* *Smilax pseudochina* *Medeola virginiana*
Bitternut Hickory	**Juglandaceae** *Carya cordiformis*
Blackjack Oak Willow Oak Water Oak Laurel Oak	**Fagaceae** *Quercus marilandica* *Quercus phellos* *Quercus nigra* *Quercus laurifolia*
Sassafras	**Lauraceae** *Sassafras albidum*
Sweetgum	**Hamamelidaceae** *Liquidambar styraciflua*
Blackberry Blackcherry	**Rosaceae** *Rubus allegheniensis* *Prunus serotina*
Poison Ivy Winged Sumac	**Anacardiaceae** *Toxicodendron radicans* *Toxicodendron copallinum*

Table 3-1
Species List - Vegetation Survey,
July 13 to 19, 1992,
Salmon Site
(Page 2 of 3)

Common Name	Scientific Name
Red Maple	**Aceraceae** *Acer rubrum*
Water Tupelo Black Gum Flowering Dogwood	**Cornaceae** *Nyssa aquatica* *Nyssa sylvatica* *Cornus florida*
White Ash	**Oleaceae** *Fraxinus americana*
Boneset Tickseed Sunflower Ragweed Goldenrod Bearded Tickseed	**Compositae** *Eupatorium perfoliatum* *Bidens* sp. *Ambrosia artemisiifolia* *Solidago* sp. *Bidens aristosa*
Buttonbush	**Rubiaceae** *Cephalanthus occidentalis*
Pawpaw	**Annonaceae** *Asimina triloba*
Baldcypress	**Taxodiaceae** *Taxodium distichum*
Saw Palmetto	**Palmae** *Serenoa repens*
Common Persimmon	**Ebenaceae** *Diospyros virginiana*
Yellow Poplar Southern Magnolia	**Magnoliaceae** *Liriodendron tulipifera* *Magnolia grandiflora*
Southern Bayberry	**Myricaceae** *Myrica cerifera*
American Holly	**Aquifoliaceae** *Ilex opaca*
Loblolly Pine Longleaf Pine Shortleaf Pine	**Pinaceae** *Pinus taeda* *Pinus palustris* *Pinus echinata*
Hooded Pitcher-Plant	**Ericaceae** *Sarracenia minor*

Table 3-1
Species List - Vegetation Survey,
July 13 to 19, 1992,
Salmon Site
(Page 3 of 3)

Common Name	Scientific Name
Common Cattail	**Typhaceae** *Typha latifolia*
Common Pipewort	**Eriocaulaceae** *Eriocaulon septangulare*
Arrow-Arum Jack-in-the-Pulpit	**Araceae** *Peltandra virginica* *Arisaema triphyllum*
Snowy White Orchid	**Orchidaceae** *Habenaria nivea*
Rattlebox Partridge Pea	**Leguminosae** *Crotalaria sagittalis* *Cassia fasciculata*
Box Huckleberry	**Ericacea** *Gaylussacia brachvcera*
Marsh Pink	**Gentianaceae** *Sabatia stellaris*
Japanese Honeysuckle	**Caprifoliaceae** *Lonicera japonica*
Common St. Johnswort	**Guttiferae** *Hypericum perforatum*
Old Field Toadflax	**Scrophulariaceae** *Linaria canadensis*
White Baneberry	**Ranunculaceae** *Actaea pachypoda*
Green Amaranth	**Amaranthaceae** *Amaranthus retroflexus*

Table 3-2
Species List - Avian Survey,
July 13 to 19, 1992,
Salmon Site
(Page 1 of 2)

Common Name	Scientific Name
Wood Duck	*Aix sponsa*
White Ibis	*Eudocimus albus*
Cattle Egret	*Bubulcus ibis*
Green Heron	*Butorides striatus*
Broad-Winged Hawk	*Buteo platypterus*
Red-Tailed Hawk	*Buteo jamaicensis*
Mississippi Kite	*Ictinia mississippiensis*
Barred Owl	*Strix varia*
Black Vulture	*Coragypes atratus*
Northern Bobwhite	*Colinus virginianus*
Wild Turkey	*Meleagris gallopavo*
Common Flicker	*Colaptes auratus*
Red-Bellied Woodpecker	*Melanerpes carolinus*
Downy Woodpecker	*Picoides pubescens*
Great Crested Flycatcher	*Myiarchus crinitus*
Eastern Wood Pewee	*Contopus virens*
Northern Rough-Winged Swallow	*Stelgidopteryx serripennis*
Chimney Swift	*Chaetura pelagica*
Common Nighthawk	*Chordeiles minor*
Bluejay	*Cyanocitta cristata*
American Crow	*Corvus brachyrhynchos*
Carolina Chickadee	*Parus carolinensis*
Tufted Titmouse	*Parus bicolor*
Carolina Wren	*Thryothorus ludovicianus*
Northern Mockingbird	*Mimus polyglottos*
Brown Thrasher	*Toxostoma rufum*
American Robin	*Turdus migratorius*
Wood Thrush	*Hylocichla mustelina*
Blue-Gray Gnatcather	*Polioptila caerulea*
Pine Warbler	*Dendroica pinus*
Common Yellow-Throat	*Geothlypis trichas*
Hooded Warbler	*Wilsonia citrina*
Swainson's Warbler	*Limnothlypis swainsonii*
Prairie Warbler	*Dendroica discolor*

Table 3-2
Species List - Avian Survey,
July 13 - 19, 1992,
Salmon Site
(Page 2 of 2)

Common Name	Scientific Name
Scarlet Tanager	*Piranga olivacea*
Summer Tanager	*Piranga rubra*
Northern Cardinal	*Cardinalis cardinalis*
Blue Grosbeak	*Guiraca caerulea*
Indigo Bunting	*Passerina cyanea*
House Finch	*Carpodacus mexicanus*
Rufous-Sided Towhee	*Pipilo erythrophthalmus*
Bachman's Sparrow	*Aimophila aestivalis*

Table 3-3
Species List - Reptiles and Amphibian Survey,
July 13 to 19, 1992,
Salmon Site

Common Name	Scientific Name
Amphibians	
Eastern Narrow-Mouthed Toad	*Gastrophyrne carolinensis*
Squirrel Treefrog	*Hyla squirella*
Slimy Salamander	*Plethodon glutinosus*
Alabama Waterdog	*Necturus alabamensis*
Bronze Frog	*Rana clamitans clamitans*
River Frog	*Rana heckscheri*
Southern Dusky Salamander	*Desmognathus auriculatus*
Southern Cricket Frog	*Acris gryllus gryllus*
Four-Toed Salamander	*Hemidactylium scutatum*
Reptiles	
Southeastern Five-Lined Skink	*Eumeces inexpectatus*
Gopher Tortoise	*Gopherus polyphemus*
Green Anole	*Anolis carolinensis*
Ground Skink	*Scincella laterale*
Fence Lizard	*Sceloporus undulatus*
Eastern Box Turtle	*Terrapene carolina carolina*
Northern Brown Snake	*Storeria dekayi dekayi*
Water Snake (unidentified)	*Nerodea sp.*
Western Cottonmouth	*Agkistrodon piscivorus leucostoma*
Southern Toad	*Bufo terrestris*

Table 3-4
Species List - Mammalian Survey,

Table 3-4
Species List - Mammalian Survey,
July 13 to 19, 1992,
Salmon Site

Common Name	Scientific Name
Armadillo[abc]	*Dasypus novemcinctus*
Big Brown Bat[a]	*Eptesicus fuscus*
Myotis[a]	*Myotis* sp.
Eastern Cottontail[abc]	*Sylvilagus floridanus*
Squirrel[ad] (Grey or Fox)	*Sciurus* sp.
Coyote[b]	*Canis latrans*
Domestic Dog[b]	*Canis familiaris*
Domestic Cat[b]	*Felis catus*
White-Tailed Deer[abc]	*Odocoileus virginianus*

[a]Identified by sighting
[b]Identified by tracks
[c]Identified by scat
[d]Identified by sound

Table 3-5
Species Collected,
February 22 to 28, 1993,
Salmon Site

Common Name	Scientific Name
White-Tailed Deer	*Odocoileus virginianus*
Raccoon	*Procyon lotor*
Opossum	*Didelphis marsupialis*
Eastern Cottontail	*Sylviagus floridanus*
Marsh Rabbit	*Sylviagus palustris*
Pine Vole	*Pitymus pinetorum*
Short-Tailed Shrew	*Blarina brevicauda*
Eastern Woodrat	*Neotoma floridana*
Mice	*Peromyscus* sp.

Table 3-6
Fish Biosurvey Results,
July 13 to 19, 1992,
Salmon Site

Common Name	Genus Species	Aquatic Survey Stations												
		BeP-1	BeP-2	GRC-1	GRC-2	HHC-1	HHC-2	HOP-1	GtP-1	HMC-1	HMC-2	HMC-3	HMC-4	HMC-5
Bass	Family: Centrarchidae	X												
Black Banded Darter	*Percina nigrofasciata*			X										
Blackstripe Topminnow	*Fundulus notatus*						X[a]							
Blacktail Shiner	*Notropis venusta*												X	
Bluegill	*Lepomis macrochirus*			X				X[a]						
Broadstripe Topminnow	*Fundulus euryzonus*												X	
Brown Bullhead	*Ameiurus nebulosus*			X								X		
Cherryfin Shiner	*Lythrurus roseipinnis*			X[a]	X[a]		X				X[a]	X[a]	X[a]	X[a]
Chestnut Lamprey	*Ichthyomyzon castaneus*				X								X	X
Flagfin Shiner	*Notropis signipinnis*			X						X[a]	X	X		
Gulf Darter	*Etheostoma swaini*											X		
Mosquito Fish	*Gambusia affinis*	X[a]	X											
Silverstripe Shiner	*Notropis stilbuis*	X											X	
Starhead Topminnow	*Fundulus notti*	X	X[a]										X	
Yellow Bullhead	*Ameiurus natalis*												X	
Alabama Waterdog	*Necturus alabamensis*												X	
Lesser Siren	*Siren intermedia*											X		

[a]Indicates dominant species at that station

3.1.3 Species of Special Concern

Only two species listed as either threatened, endangered, or sensitive are known to inhabit the SS. These are the gopher tortoise (*Gopherus polyphemus*), (federally-listed threatened and state-listed endangered) and Bachman's sparrow (federal Category 2, information insufficient to list). However, the habitat supporting these species on the SS is not unique in any way, but is part of an expanse of similar pine forest that is common throughout the southeastern part of the state. A complete listing of the survey methods and results for threatened and/or endangered species is presented in the Results of a Preliminary Survey for Threatened and Endangered Species and Floodplains/Wetlands at the Tatum Dome Test Site Lamar County, Mississippi (IT, 1992a).

3.2 Constituents of Potential Concern Selection

A multistep method was used to identify COPC for the site because of the large quantity of analytical data available for both chemical and radiological constituents.

The first step involved identifying a set of reported concentrations that were of acceptable quality for use in the environmental evaluation. This was accomplished through (a) evaluations of analytical methods, (b) evaluations of the quality of the data with respect to detection limits, and (c) laboratory and validation qualifiers and blanks.

The second step was to reduce the data set to those constituents that frequently had higher concentrations in the on-site samples than were found in the reference sample. This is an indication that these constituents are present due to site-related activities.

The third step involved identifying COPC within the soil, sediment, and surface water, which were of greatest concern with respect to bioaccumulation in ecological receptors based on the octanol/water partition coefficient (K_{ow}).

The list of COPC was further reduced, on the basis of water or sediment quality criteria and detection in biota tissue samples, to a final list of contaminants of concern. A COC was selected if water or sediment quality criteria, which are developed to be protective of all aquatic life, were exceeded, or if the constiuent was detected in tissues. If no aquatic criteria existed for water or sediment COPC, but the constituent was present in fish tissue, it was considered a COC. Since no mandated soil quality criteria exist, and COPC were detected in animal and plant tissues, these were carried through as COC. This list of COC was then used to assess contaminant-related risk at the SS.

3.2.1 Data Collection

3.2.1.1 Reference Samples

Reference sites for surface water and sediment were established on Half Moon Creek (HMC-1), located at the southern boundary; Grantham Creek (GrC-1), located at the eastern boundary; and Hickory Hollow Creek (HHC-1), located at the western boundary of the SS. An unnamed catfish pond located off site (ReP-1) was used as the source of the northern reference sample. Figure 3-1 shows the locations of the surface-water sampling stations. The field adjacent to the catfish pond was used as the reference location for surficial soils. Figure 3-2 shows the location of the soil sampling stations at the SS. Reference sampling was conducted concurrently with the on-site sampling events.

3.2.1.2 Sediment and Surface Water

Sediment samples were collected during the period of February 22 to 28, 1993. Only 15 of the 22 stations originally identified in the Environmental Evaluation Work Plan (EEW) were sampled during the investigation (Table 3-7). During the initial site reconnaissance, it was determined that the ponds in SA-2 and SA-4 no longer existed. Swamps identified in previous investigations were also absent or were now part of the Beaver Pond. The large size of the Beaver Pond required that three sampling stations be used instead of one. Station BeP-1 was located on the northern end of the pond near the E-6 Decontamination Pad; station BeP-1a at the southeastern end near the Bleed Down Plant Area and E-14 Pad; and station BeP-2 was located in a small pond connected to the main pond by a galvanized steel culvert beneath Half Moon Circle Road. This portion of the Beaver Pond is located near the E-14 Pad and Mud Pits. Sediment samples were analyzed for organic, inorganic, and radiological constituents. In addition, grain size and total organic carbon were measured. Figure 3-1 shows the location of these sampling stations.

Surface water samples were analyzed for organic, inorganic, and radiological constituents, as well as general water-quality chemistry parameters. The biochemical oxygen-demand (BOD) analyses were conducted after holding time had expired. The original BOD data were supplemented with five additional samples collected on April 28, 1993. These five samples were collected from HMC-4, HOP-1, HHC-2, HMC-2, and GrC-1. Figure 3-1 shows the locations of these sampling stations. A summary of the types of samples collected for the aquatic investigation is provided in Table 3-8.

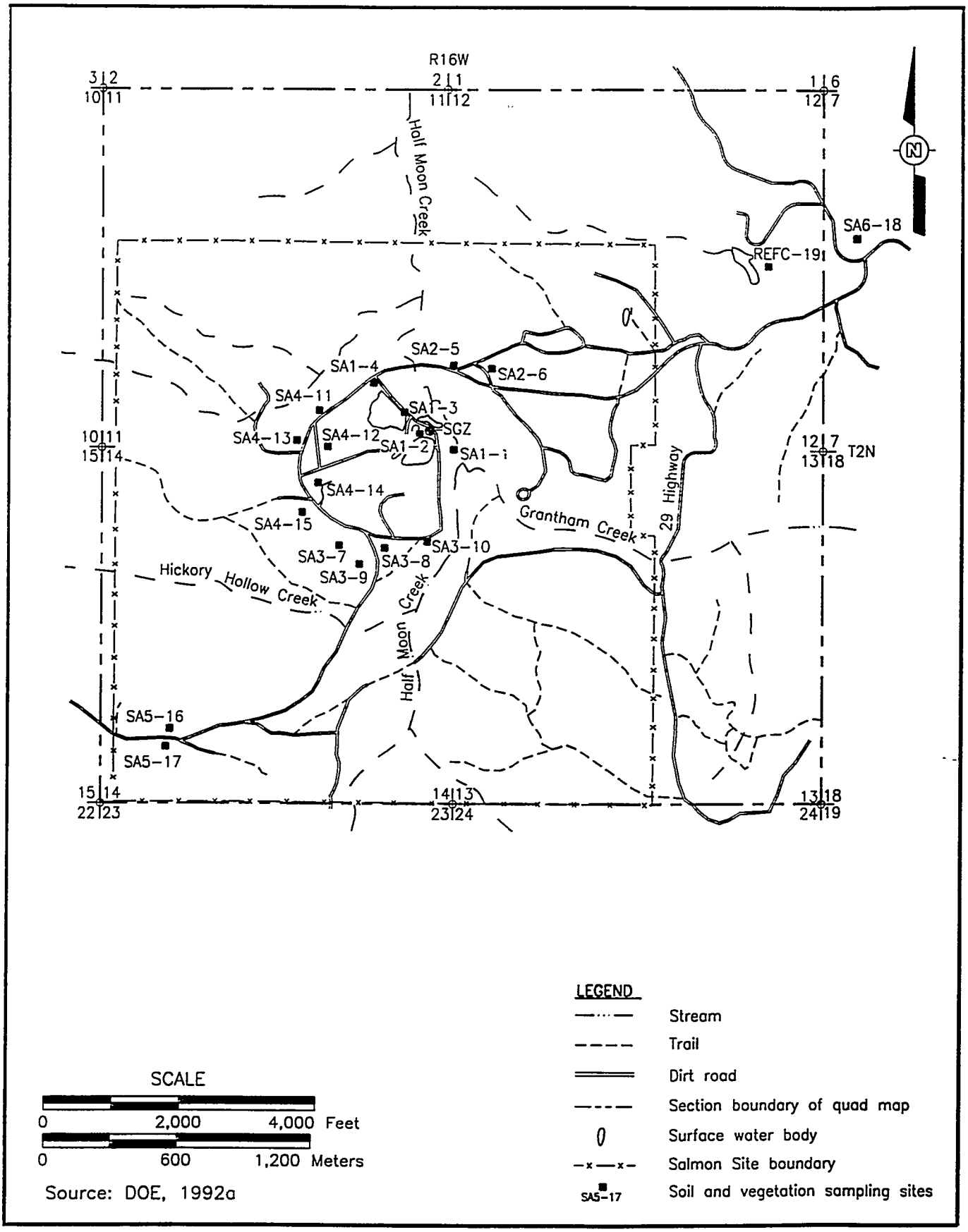

Figure 3-2
Salmon Site Soil- and Vegetation-Sampling Locations

Table 3-7
Surface Water and Sediment Sample Locations for the Aquatic Ecosystem Investigation, Salmon Site

HMC-1	Half Moon Creek just within the southern boundary of the Salmon Site (SS) (Reference).
HMC-2	Half Moon Creek approximately ½ mile upstream of the confluence of Hickory Hollow Creek and Half Moon Creek. This station is downstream of surficial groundwater flow from Source Area (SA)-5.
HMC-3	Half Moon Creek approximately 1/8 mile upstream of the confluence of Grantham Creek and Half Moon Creek. This station is in an area of surficial groundwater flow from SA-5.
HMC-4	Half Moon Creek approximately ½ mile downstream of the confluence of Grantham Creek and Half Moon Creek at the bridge where the Main Road crosses Half Moon Creek. This station is in an area just downstream of surficial groundwater flow from SA-1.
HMC-5	Half Moon Creek just within the northern boundary of SS.
HHC-1	Hickory Hollow Creek just within the western boundary of SS (Reference).
HHC-2	Hickory Hollow Creek approximately ½ mile upstream of the confluence of Hickory Hollow Creek and Half Moon Creek. This station is downstream of surficial groundwater flow from SA-5.
GrC-1	Grantham Creek just within the eastern boundary of SS (Reference).
GrC-2	Grantham Creek approximately 1/8 mile upstream of the confluence of Grantham Creek and Half Moon Creek. This station is an area of surficial groundwater flow from SA-2.
BeP-1	Beaver Pond 1 on the northern end of the Pond. This pond receives runoff and groundwater discharge from potential "hot spots" within SA-1 and SA-2.
BeP-1a	Beaver Pond location 2 on the southern end of the Pond.
BeP-2	Beaver Pond location 3 on the small pond connected to the Beaver Pond.
HOP-1	Half Moon Creek Overflow Pond. This pond has received runoff and groundwater discharge that was contaminated with tritium from surface ground zero.
ReP-1	Reference Pond. This privately-owned pond is located off the northeast corner of the SS (Reference).

Table 3-8
Summary of Samples Collected for the Aquatic Ecosystem Investigation, Salmon Site

Salmon Site Surface Water Station	Water Quality Parameters	Analytical Samples	Grain Size and Sediment Analysis	Benthic Ecology Analysis	Fish Tissue Analysis	Toxicity Tests
HMC-1	X[a]	X	X	X	X[b]	X
HMC-2	X	X	X	X	X[b]	X
HMC-3	X	X	X	X	X[b]	X
HMC-4	X	X	X	X	X[b]	X
HMC-5	X	X	X	X	X[b]	X
HHC-1	X	X	X	X	NA[c]	X
HHC-2	X	X	X	X	NA	X
GrC-1	X	X	X	X	X[b]	X
GrC-2	X	X	X	X	X[b]	X
BeP-1	X	X	X	X	X	X
BeP-1A	X	X	X	X	X	X
BeP-2	X	X	X	X	X	X
GtP-1	X	X	X	X	NA	X
HOP-1	X	X	X	X	NA	X
Ref. Pond	X	X	X	X	X	X

[a]Sample collected
[b]Fish tissue sample represents a composite sample of the organisms collected for that creek or pond system.
[c]No sample was obtained from the collection effort.

3.2.1.3 Surficial Soil

Surficial soil samples were collected during the period of April 26 to 30, 1993, and analyzed for chemical and radiological constituents. All 18 sites (excluding ponds, mud pits, and waste burial pits) located on site were sampled individually and not composited as originally planned. The reference site was also sampled. The 19 soil sample stations are listed in Table 3-9.

Five soil samples were collected at each site: one at the approximate center of site activity and one each 30 m (100 ft) from the center to the north, south, east, and west. The sample designator A was given to the north sample, B to the west sample, C to the center sample, D to the east sample, and E to the south sample.

Soil samples were collected from the top 10 centimeters (cm) (4 inch [in.]) of soil, excluding vegetation and leaf litter. At each location, the leaf litter was removed by gently pushing it aside. After an area of approximately 10-cm diameter had been cleared, a stainless-steel shovel was used to excavate a 10-cm-deep hole. This hole was then tested with an organic vapor analyzer before sampling. After being deemed safe from organic vapors in the breathing zone, sections of the undisturbed wall were removed and placed in the sample containers. Samples were then shipped on ice to the International Technology Analytical Services (ITAS) Laboratory for analysis. Table 3-10 summarizes the data collected for the terrestrial investigation.

3.2.2 Data Evaluation

3.2.2.1 Analytical Methods

The data sets used in the quantitative evaluations have been developed according to a standard set of sensitive, chemical-specific methods. The standard methods apply both to the collection of samples and the analyses thereof. Due to the lack of information on the methods used to collect and analyze the earlier samples, only the most recent data were used in the quantitative evaluations. The data from earlier studies were only used to guide the development of the EEW.

Analytical results that are not specific for a particular compound or that are the result of insensitive analytical methods are useful for qualitative assessments but are generally not appropriate for quantitative evaluations. Likewise, results of analytical methods associated

Table 3-9
Sample Locations for the Terrestrial Ecosystem Investigation, Salmon Site

Source Area (SA)-1 - Surface Ground Zero	
SA-1-1	Station 1A Surface Ground Zero
SA-1-2	Bleed Down Plant Area
SA-1-3	East Substation
SA-1-4	E-6 Decon Area
SA-2 - Northern Disposal Area	
SA-2-5	Clean Burn Pit
SA-2-6	Gas Pump
SA-3 - Southern Disposal Area	
SA-3-7	Big Chief Drilling Storage
SA-3-8	Site E-2 and E-7
SA-3-9	Government Storage Area 2 (Drilling Storage Yard)
SA-3-10	Government Storage Area 3
SA-4 - Western Disposal Area	
SA-4-11	CH Fuel Storage
SA-4-12	Cable Storage
SA-4-13	North Substation
SA-4-14	South Substation
SA-4-15	West Substation
SA-5 - Injection Well Area	
SA-5-16	HT-2 Well Area
SA-5-17	HT-2M Well Area
SA-6 - Helicopter Pad and Storage Area	
SA-6-18	Helicopter Pad and Storage Area
Reference Area	
Ref-19	Reference Station

Table 3-10
Summary of Samples Collected for the Terrestrial Ecosystem Investigation, Salmon Site

Salmon Site	Soil Analyses	Field Biota Observations	Vegetation Tissue Analyses	Mammalian Tissue Analyses	Toxicity Tests
Source Area (SA) 1-1	X[a]	X	X	X[b]	X[a]
SA-1-2	X[a]	X	X	X[b]	X[a]
SA-1-3	X[a]	X	X	X[b]	X[a]
SA-1-4	X[a]	X	X	X[b]	X[a]
SA-2-5	X[a]	X	X	X[b]	X[a]
SA-2-6	X[a]	X	X	X[b]	X[a]
SA-3-7	X[a]	X	X	X[b]	X[a]
SA-3-8	X[a]	X	X	X[b]	X[a]
SA-3-9	X[a]	X	X	X[b]	X[a]
SA-3-10	X[a]	X	X	X[b]	X[a]
SA-4-11	X[a]	X	X	X[b]	X[a]
SA-4-12	X[a]	X	X	X[b]	X[a]
SA-4-13	X[a]	X	X	X[b]	X[a]
SA-4-14	X[a]	X	X	X[b]	X[a]
SA-4-15	X[a]	X	X	X[b]	X[a]
SA-5-16	X[a]	X	X	X[b]	X[a]
SA-5-17	X[a]	X	X	X[b]	X[a]
SA-6-18	X[a]	X	X	X[b]	X[a]
Ref-19	X[a]	X	X	X[b]	X[a]

[a]Analyses or tests conducted on five replicates from the site.
[b]Collected as site-wide samples.

with unknown, few, or no Quality Assurance/Quality Control procedures were eliminated from quantitative consideration.

3.2.2.2 Detection Limits

All analyzed chemicals were evaluated with respect to their quantitation or detection limits. This evaluation was either conducted by the laboratory or later during data validation.

Constituents that were analyzed for, but not detected, were eliminated from the COPC list; however, before eliminating the constituents from further evaluation, several points were considered. Sample quantitation limits of known site-related constituents may be greater than corresponding standards, criteria, or other levels of concern. Therefore, these constituents may be present at concentrations greater than the corresponding reference concentrations. A sample-specific detection limit may be significantly higher than positively detected values in other samples in a data set. In other words, only a limited number of samples may test positive for a particular constituent.

When determining exposure concentrations at the SS, positively detected results were considered together with nondetected results. If there was reason to believe that the constituent was present in a sample at a concentration below the detection limit, then half the detection limit was used as the estimated concentration for that sample (EPA, 1989a). This method assumes that constituent concentrations are uniformally distributed in the range of concentrations below the detection limit. Therefore, half the detection limit is the arithmetic average of all such concentrations and is also the maximum likelihood estimator.

If the site-specific information indicated that a constituent was not likely to be present in a sample, a value of zero may have been used as the estimated concentration. In cases when the constituents were anthropogenic in nature and site-specific information indicated localized usage and limited transport from areas of use, the use of zero as the estimated concentration would be considered acceptable in areas of no known usage of these constituents.

Those constituents that were not detected in any samples of a particular medium were eliminated from further consideration in the quantitative evaluations. The outcome of these steps was a data set that contained only constituents which were detected in at least one sample from each medium.

3.2.2.3 *Qualified Data*

Qualifiers may be attached to the analytical results by the laboratory conducting the analyses or by persons performing the data validation. These qualifiers generally indicate questions concerning constituent identity, concentration, or both.

Because the data validation process assesses the effect of quality control issues on data usability, validation data qualifiers supersede the laboratory qualifiers. When an analytical result has laboratory and validation qualifiers that appear contradictory, only the validation qualifier is considered. Laboratory qualifiers are valid unless they have been superseded during data validation.

The qualifiers used in the SS data sets are presented in the following paragraphs. Organic, inorganic, and radiochemical qualifiers are presented separately since some of the qualifiers have different meanings for different data sets.

The organic qualifiers are as follows:

- B - indicates that the constituent was found in an associated blank, as well as in the sample. Use of these data is discussed in the next subsection.
- E - indicates that the concentration exceeded the calibration range of the Gas Chromatograph/Multispectral Scanner. The use of this qualifier implies uncertainty in concentration but not in identity. These concentrations were treated as positive detections.
- J - indicates that the reported results were quantitatively estimated. These concentrations were treated as positive detections.
- NJ - indicates that there was presumptive evidence that a Tentatively Identified Compound (TIC) was present at an estimated concentration. The use of this data is discussed in a subsequent subsection.
- NP - indicates that data validation was not performed. Laboratory qualifiers were valid for these data.
- R - indicates that the data were rejected. These data were unusable and could not be used in the quantitative evaluations.
- U - indicates that the constituent was analyzed for but not detected. The use of these data was discussed in the previous subsection.
- UJ - indicates that the reported quantitation limit was quantitatively estimated. These data were treated as not detected and were used as indicated in the previous subsection.

- 5U or 10U - indicates that the 5-times and 10-times rule (EPA, 1989a) was applied due to blank contamination. Use of these data is discussed in the following subsection.

The inorganic qualifiers are as follows:

- B - indicates that the reported value was below the Contract Required Detection Limit (CRDL), but above the Instrument Detection Limit. The use of this qualifier indicated uncertainty in concentration but not in identity. These data were treated as positive detections.

- E - indicates that the Inductively Coupled Plasma dilution (%D) was out of control limits. This qualifier indicated uncertainty in the reported concentration. These data were treated as positive detections.

- J - indicates that the reported result was qualitatively estimated. These data were treated as positive detections.

- N - indicates that the spiked sample recovery was not within control limits. This qualifier indicated uncertainty in the reported concentration. These data were treated as positive detections.

- NJ - indicates that there was presumptive evidence that a TIC was present at an estimated concentration.

- NP - indicates that data validation was not performed. Laboratory qualifiers were valid for these data.

- R - indicates that the data were rejected. These data were unusable and could not be used in the quantitative evaluations.

- S - indicates that the reported value was determined by the Method of Standard Addition. No uncertainty in either concentration or identity was implied by the use of this qualifier. Data were treated as positively detected.

- U - indicates that the constituent was analyzed for but not detected.

- UJ - indicates that the reported quantitation limit was quantitatively estimated. These data were treated as not detected and were used as indicated in the previous subsection.

- 5U or 10U - indicates that the 5-times and 10-times rule was applied due to blank contamination. Use of these data is discussed in the following subsection.

- W - indicates that the postdigestion spike for furnace atomic absorption analysis was out of control limits. This qualifier indicated uncertainty in the reported concentration. These data were treated as positive detections.

The radiochemical constituent qualifiers are as follows:

- J - indicates that the reported result was qualitatively estimated. These data were treated as positive detects.
- ND - indicates that the constituent was not detected since the result was below the Minimum Detectable Activity (MDA). Such data were treated as not detected. Appropriate estimated concentrations were used as discussed in the previous subsection.
- NP - indicates that data validation was not performed. Laboratory qualifiers were valid for these data.
- U - indicates that the compound was not detected at the specified detection limit. Appropriate estimated concentrations were used as discussed in the previous subsection.
- UJ - indicates that the reported quantitation limit was quantitatively estimated. These data were treated as not detected and were used as indicated in the previous subsection.

3.2.2.4 Blanks

Blank samples provide a means to check for contamination that may have been introduced into a sample in the field while the samples were being collected and transported or while in the laboratory during sample preparation or analysis. The concentrations of constituents detected in blanks must be compared with the concentrations of the same constituents detected in site samples in order to prevent the inclusion of nonsite-related constituents into the quantitative evaluation of COPC.

As part of the data validation process, an analyte was not considered further if the maximum sample concentration did not exceed 10 times the highest blank for all common laboratory contaminants (acetone, 2-butanone, methylene chloride, toluene, and phthalates) or 5 times the highest blank for other chemicals (EPA, 1989a).

3.2.2.5 Frequency of Detection

Constituents that are infrequently detected may be artifacts in the data due to sampling or analytical procedures or from other means of contamination. Chemicals that were detected in 5 percent or less of the on-site samples for a particular medium were eliminated as COPC. Less than 20 surface water or sediment samples were taken and analyzed. In these cases, a constituent was considered a COPC unless the detected concentration, when averaged with half the detection limit for nondetected samples, produced a 95-percent upper confidence limit (UCL) that was within the reference range or less than the actual detection limit for the reference samples or that could be eliminated on another basis as discussed below.

3.2.2.6 *Essential Nutrients*

Calcium, sodium, potassium, iron, and magnesium aré generally eliminated as constituents of potential ecological concern because they are essential elements for animal and plant life and are only toxic at very high concentrations, as supported by EPA (1989a). However, these may be products of site activities, specifically abandoned mud pits from rotary drilling. Further examination of these elements indicates that calcium was the only analyte among the five mentioned above for which the 95-percent upper confidence limit of the mean for site samples in surficial soils exceeded concentrations in soil samples collected from reference locations.

Limited data exist on the toxicity of calcium. As indicated by the National Research Council (1980), toxicity of calcium is dependent on the level of dietary potassium, and on the age and reproductive status of the animal. Maximum tolerance levels (MTL) for calcium are 2 percent of the diet for cattle, sheep, horse, and rabbits; 4 percent for swine; and 1 percent of the diet for poultry (National Research Council, 1980). If a value of 2 percent is used for the white-tailed deer (body weight of 90.72 kilogram [kg]) (Harestad and Bunnel, 1979); estimated ingestion rate of 2,319 grams per day (g/day) (EPA, 1993) and 1 percent for the red-tailed hawk (body weight of 1.126 kg [Dunning, 1994]; ingestion rate of 58.9 g/day [EPA, 1993]), the MTLs for these receptorsare estimated to be 46.4 grams (g) of calcium per day and 0.707 g of calcium per day for the deer and hawk, respectively. Further, assuming that the entire diet of the deer consists of plant matter and using a soil-to-plant transfer factor of 3.5 (Baes et al., 1984), the deer would be exposed to 6.54 g of calcium per day. Calcium is, therefore, not expected to be toxic to this herbivorous mammal. If it assumed that the hawk consumes only herbivorous mice from the site (Peromyscus maniculatus, body weight of 21 g [Millar, 1989 as cited in EPA, 1993]; ingestion rate of 3.46 g/day [EPA, 1993]; absorption coefficient of 0.80 assumed [Ganong, 1979]), the hawk would be exposed to 0.022 g of calcium per day. As with the deer, calcium at the site is not expected to be toxic to the top avian receptor. This provides quantitative support to the original contention that essential nutrients be eliminated as COPC because they are generally toxic only at very high concentrations.

With reference to sediments, calcium, magnesium and sodium, 95-percent UCL of the mean of site samples exceeded reference ranges. For surface water, site iron, calcium, magnesium and sodium exceeded the reference ranges. Potassium was not detected at the CRDL in water or sediment. No federally mandated Ambient Water Quality Criteria (AWQC) or sediment quality criteria exist for calcium, sodium or magnesium. In general, these elements are not considered toxic to aquatic life, a contention supported by the fact that no toxicity benchmarks have been established or mandated by EPA. One location had particularly high levels of calcium, sodium

and magnesium in surface water (Half Moon Creek overflow pond), and another (Beaver Pond) also had some high values in surface water and sediments. These two areas are likely to have received materials from nearby rotary drilling mud pits. Since evaluation of toxicity of calcium, sodium and magnesium in aquatic environments has not been promulgated by EPA, it will be assumed that in general, the aquatic environments at the SS do not have sufficiently high concentrations to cause toxicity, and will not be carried through as COPC. The Beaver Pond and Half Moon Creek overflow pond, however, should be held under consideration as "hot spots" for these constituents.

An AWQC of 1,000 ug/ℓ does, however, exist for the chronic exposure of freshwater biota to iron. The 95-percent UCL of the mean for site surface water samples exceeded the reference range and the AWQC, and iron was therefore carried through the assessment as a COPC in surface water.

3.2.2.7 Water Chemistry

General water chemistry parameters, such as sulfates, carbonate, bicarbonate, chloride, dissolved solids, and fluoride, are also generally toxic only at very high levels. Measurements of these are useful in determining the nature of a water body being studied, but are not considered COPCs for the SS.

3.2.2.8 Physiochemical Properties and Bioaccumulation Potential

The K_{ow} provides a useful screening tool for organic compounds on the basis of the bioaccumulation potential. In general, a higher K_{ow} relates to a higher bioaccumulation potential (i.e., a tendency to occur in biota at higher concentrations than the surrounding environment). Garten and Trabalka (1983) state that compounds with log K_{ow} values of less than 3.5 generally do not bioaccumulate in the fat tissues of mammals and birds by a factor of greater than 0.3. In addition, according to Thomann (1989), bioaccumulation in aquatic food chains is not significant for compounds with log K_{ow} values of up to 5. Finally, EPA (1993) has correlated food chain multipliers with log K_{ow} values and trophic level in an aquatic system. Compounds with log K_{ow} values of less than or equal to 3.9 are given food chain multipliers of 1 or less for each of the three trophic levels, which include zooplankton in level 2, small fish in level 3, and piscivorous fish in level 4. It is also recommended that trophic levels 3 and 4 be used for wildlife, depending on the specific bird or mammal in question. For these reasons, the conservative log K_{ow} value of 3.5 was used to blanket all potential bioaccumulation pathways. Constituents with a log K_{ow} less than 3.5 were not therefore, considered as COPC via bioaccumulation pathways.

3.2.2.9 Reference Data

Reference data were gathered during the BERA investigation using methods and procedures identical to those used for gathering site samples. Reference data were collected from areas not influenced by the site. They are collected from each medium of concern (soil, sediment, and surface water) from areas that could not have received contamination from the site, but that do have the same basic characteristics as the medium of concern on site.

Reference soil samples were collected from an area off site and upgradient of the site. Reference surface water and sediment samples were collected from each of the three on-site creeks (Hickory Hollow Creek, Half-Moon Creek, and Grantham Creek) at points where the creeks enter the site and from a reference pond located off site and upgradient of the site.

Site data were evaluated with respect to reference data by comparing the calculated exposure point concentration (EPC) for each constituent to the range of reference and background concentrations for that constituent. The EPC is a conservative estimate of the most probable concentration in a particular media to which biota may be exposed. It is calculated as the 95-percent UCL on the arithmetic mean of the positive detections and half the detection limit for samples in which the analyte was not detected, using the following equation:

EPC = 95-percent UCL of the mean =

$$\text{mean} + [t_{0.05(2)(n-1)} \ (s/\sqrt{(n)})]$$

Where: n = number of samples
$t_{0.05(2)(n-1)}$ = value from Student's t table corresponding to $\alpha_{(2)} = 0.05$
n-1 = degrees of freedom
s = standard deviation

If the EPC is within the range of background or reference data, the constituent's presence is considered to be ambient and not site-related. The constituent is then removed from the quantitative evaluation.

A summary of the detected constituents in sediment, surface water, and surficial soil are presented in Tables 3-11 through 3-14, 3-15 through 3-16, and 3-17 through 3-20, respectively. The EPC and reference ranges are included for data evaluation comparisons. Those constituents present on site at levels greater than ambient are considered potentially site-related constituents.

3.2.3 Analytical Results

Analyte concentrations were measured in sediment, surface water, surficial soil, vegetation, and wildlife tissues. All data are presented in Appendix A of this report and summarized below. The summary tables presented in this section use the data evaluation and data estimation procedures for nondetected concentrations discussed in the preceding sections.

3.2.3.1 Sediment

Sediment data are presented in Appendix A. Tables A-3.1 through A-3.9 present the inorganic, pesticide and polychlorinated biphenyls (PCB), volatile, semivolatile, gross alpha and beta, gamma spectroscopy, tritium, grain size, and wet chemistry data sets, respectively.

The majority of inorganic constituents detected in sediments are naturally occurring soil constituents and were found at reference concentrations. Cadmium was found at the Beaver Pond in concentrations above background. Concentrations in the Beaver Pond sediments ranged from 1.7 to 9.0 milligrams per kilogram (mg/kg) cadmium. Cadmium was not detected in the Reference Pond sediments. The Beaver Pond sediments had the highest levels of many of the inorganic constituents: aluminum (13,000 mg/kg), arsenic (7.7 mg/kg), barium (181 mg/kg), calcium (5,630 mg/kg), lead (20.2 mg/kg), magnesium (1,380 mg/kg), manganese (583 mg/kg), sodium (1,230 mg/kg) and zinc (111 mg/kg). Chromium was highest in the Half Moon Creek Overflow Pond sediment (14.8 mg/kg). Concentrations of inorganics in sediments on the SS are summarized in Table 3-10.

No pesticides or PCBs were detected in sediments on the SS. However, 3 organic compounds were detected in the Beaver Pond sediment samples: 2-butanone (170 micrograms per kilogram [µg/kg]), toluene (220 µg/kg) and benzoic acid (550 µg/kg, estimated concentration below the detection limit). Sediment volatile and semivolatile organic compound concentrations are summarized in Table 3-11.

Gross alpha levels above the MDA were found in sediments of the Beaver Pond (17.21 picocuries per gram [pCi/g]) and Gator Pond (22.48 pCi/g). Gross beta levels above the MDA were detected at all sediment stations (8.84 pCi/g to 23.69 pCi/g); the highest levels were found in the Gator Pond (23.69 pCi/g) and Beaver Pond (18.6 pCi/g) sediments.

Table 3-11
Summary of Inorganic Concentrations in Sediment, Salmon Site

Analyte	Frequency of Detection	Minimum Detection Limit[a] (mg/kg)[b]	Range of Detected Concentrations (mg/kg)	Reference Range[c,d] (mg/kg)	Average Concentration[d] (mg/kg)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (mg/kg)	Constituent of Potential Concern
Aluminum	12/12	--	84.40 - 13,000	195 - 6480	3,218	3,973	5,742	No[e]
Arsenic	7/12	0.30	0.30 - 7.70	0.0007(ND)[f] - 2.10	1.54	2.27	2.98	Yes
Barium	12/12	--	2.40 - 181	1.4(ND) - 19.90	42.71	54.82	77.54	Yes
Cadmium	2/12	0.25	1.70 - 9.00	0.37(ND) - 0.48(ND)	1.43	2.44	2.98	Yes
Calcium	12/12	--	179 - 5,630	190 - 451	1,195	1,682	2,264	No[g]
Chromium	9/12	0.90	1.20 - 14.80	0.45(ND) - 9.30	4.88	5.13	8 14	No[e]
Iron	12/12	--	125 - 13,000	102 - 19,300	3,564	4,692	6,545	No[d,g]
Lead	12/12	--	0.61 - 20.20	0.0034 - 5.00	6.23	6.38	10.23	Yes
Magnesium	12/12	--	39.80 - 1,380	47.20 - 152	273	378	514	No[g]
Manganese	12/12	--	6.30 - 583	0.80(ND) - 85.10	109	161	211	Yes
Sodium	11/12	60.7	59.80 - 1,230	55.40 - 80.80	238	8347	459	No[g]
Vanadium	8/12	0.40	1.40 - 27.70	0.48(ND) - 43.00	5.98	8.17	11.17	No[e]
Zinc	4/12	2.20	1.60 - 111	1.1(ND) - 2.60	15.93	31.21	35.76	Yes

[a]Minimum detection limit was selected from those samples in which the analyte was not detected.
[b]Milligram per kilogram
[c]Reference range consists of samples GrC-1, HHC-1, HMC-1, and Rep-1.
[d]Half the detection limit was used for those samples in which the analyte was not detected at the detection limit.
[e]Within reference range
[f]Not detected; value represents half the detection limit.
[g]Essential nutrient

Table 3-12
Summary of Semivolatile and Volatile Organic Concentrations in Sediment, Salmon Site

Analyte	Frequency of Detection	Minimum Detection Limit[a] (mg/kg)[b]	Range of Detected Concentrations (mg/kg)	Reference Range[c,d] (mg/kg)	Average Concentration[d] (mg/kg)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (mg/kg)	Constituent of Potential Concern
Toluene	2/15	6.00	11.00 - 220	3.50(ND)[e] - 4.50(ND)	20.03	55.44	50.73	No[d,f]
2-Butanone	2/15	13.00	5.00 - 170	6.5(ND) - 9.00(ND)	20.87			No[e]
Benzoic Acid	1/12	2,100	550	1,000(ND) - 1,150(ND)	2,300	2,684	2,858	No[d]

[a]Minimum detection limit was selected from those samples in which the analyte was not detected.
[b]Milligram per kilogram
[c]Reference concentration range consists of samples GrC-1, HMC-1, HHC-1, and ReP-1.
[d]Half the detection limit was used for those samples in which the analyte was not detected at the detection limit.
[e]Not detected; value represents half the detection limit.
[f]Log octanol/water partition coefficient (K_{ow}) less than 3.5, unlikely to bioaccumulate (Garten and Trabalka, 1983).
[g]Common laboratory contaminant

Table 3-13
Summary of Radionuclide Activity in Sediment, Salmon Site

Radionuclide	Frequency of Detection	Minimum Detection Limit[a]	Range of Detected Concentrations (pCi/g)[b]	Reference Range[c,d] (pCi/g)	Average Concentration[d] (pCi/g)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (pCi/g)	Exceeds Reference Range?
Gross Alpha	2/12	7.47	17.21 - 22.4	5.58(ND)[e] - 7.53(ND)	9.14	5.26	12.48	Yes
Gross Beta	12/12	--	8.84 - 23.69	4.3 - 12.61	12.69	5.26	16.03	Yes
Beryllium-7	1/1	--	2.40	NA[f]	NA	NA	NA	No[g]
Bismuth-212	1/1	--	0.09	0.11	NA	NA	NA	No[g]
Bismuth-214	1/1	--	0.43	NA	NA	NA	NA	No
Cadmium-109	1/2	1.10	0.70	1.00(ND) - 0.74	0.61	0.08	1.33	No[g]
Cerium-144	1/1	--	0.30	0.4(ND)	NA	NA	NA	No[g]
Cesium-137	6/12	0.01	0.07 - 1.09	0.01(ND) - 0.11(ND)	0.19	0.33	0.40	Yes
Lead-210	6/7	1.41	0.20 - 9.60	0.20 - 0.99	2.10	3.35	5.19	Yes
Lead-212	4/10	0.02	0.43 - 1.05	0.01(ND) - 1.42	0.80	0.26	1.00	No[g]
Radium-224	3/11	0.17	0.68 - 4.14	0.065(ND) - 1.26	0.87	1.17	1.66	Yes
Radium-226	1/13	0.02	0.36	0.005(ND) - 0.15(ND)	0.03	NA	NA	Yes
Radium-228	2/11	0.03	0.80	0.015(ND) - 1.16	0.81	NA	NA	No[g]
Thallium-208	6/6	--	0.04 - 0.13	0.008(ND) - 0.39	0.09	0.04	0.13	No[g]
Tritium	1/12	0.06	0.40	0.02(ND) - 0.035(ND)	NA	NA	NA	Yes

[a] Minimum Detectable Activity (MDA) was selected from those samples in which the analyte was not detected.
[b] Picocurie per gram
[c] Reference range consists of samples GrC-1, HMC-1, HHC-1, and ReP-1.
[d] Half the MDA was used for those samples in which the analyte was not detected at the MDA.
[e] Not detected; value represents half the MDA.
[f] Not applicable
[g] Maximum detected sample concentration was less than or equal to reference.

Table 3-14
Grain-Size Analysis,
Salmon Site

Sample	Percent Moisture	Percent Sand (> 0.0625 mm[a])	Percent Silt (0.062 mm to 0.0039 mm)	Percent Clay (< 0.0039 mm)
BEP-1	66.5	11.1	65.7	23.2
BEP-1A	25.0	40.8	34.5	24.7
BEP-2	27.0	86.5	7.6	5.9
AP-1	79.1	42.6	33.9	23.5
GRC-1	20.6	100.0	0.0	0.0
GRC-2	23.4	100.0	0.0	0.0
HHC-1	30.6	100.0	0.0	0.0
HHC-2	18.7	96.3	1.8	1.9
HMC-1	21.7	95.2	2.0	2.8
HMC-2	29.3	93.2	3.8	3.0
HMC-3	24.3	98.2	0.7	1.1
HMC-4	21.6	100.0	0.0	0.0
HMC-4D	22.7	100.0	0.0	0.0
HMC-5	33.6	96.3	1.8	1.9
HOP-1	26.2	91.6	4.9	3.6
REP-1	28.2	70.9	7.0	22.2

[a]Millimeter

Table 3-15
Summary of Inorganic Concentrations in Surface Water, Salmon Site

Analyte	Frequency of Detection	Minimum Detection Limit[a] (µg/ℓ)[b]	Range of Detected Concentrations (µg/ℓ)	Reference Range[c,d] (µg/ℓ)	Average Concentration[d] (µg/ℓ)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (µg/ℓ)	Constituent of Potential Concern
Aluminum	12/12	--	87.80 - 416	164 - 201	201	108	269	Yes
Arsenic	1/12	1.00	2.20	0.50(ND)[e]	0.64	0.49	0.95	Yes
Barium	12/12	--	18.80 - 76.70	19.90 - 90.40	29.83	16.07	40.04	No[f]
Cadmium	2/12	2.70	1.35 - 3.20	1.35(ND) - 3.80	1.64	0.68	2.07	No[f]
Calcium	12/12	--	569 - 24,600	573 - 3,070	3,454	6,767	7,754	No[g]
Cobalt	1/12	6.00	9.10	3.00(ND) - 10.10	3.76	1.89	4.96	No[f]
Iron	12/12	--	118 - 4,190	111 - 677	1,012	1,261	1,813	No[g]
Lead	2/12	1.00	1.10 - 1.20	0.50(ND) - 1.20	0.61	0.25	0.77	No[f]
Magnesium	12/12	--	405 - 2,220	363 - 812	729	496	1,044	No[g]
Manganese	12/12	--	21.50 - 576	34.40 - 213	113	153	210	No[f]
Mercury	1/12	0.10	0.20	0.05(ND) - 0.15	0.06	0.03	0.08	No[f]
Silver	2/12	3.50	3.50 - 5.30	1.75 - 14.60	2.19	1.10	2.89	No[f]
Sodium	12/12	--	1,640 - 56,400	1,220 - 2,160	8,011	15,926	18,130	No[g]
Zinc	3/12	3.00	3.30 - 7.20	2.15(ND) - 13.80	3.78	1.98	5.04	No[f]

[a] Minimum detection limit was selected from those samples which the analyte was not detected.
[b] Microgram per liter
[c] Reference range consists of samples GrC-1, HHC-1, HMC-1, and ReP-1.
[d] Half the detection limit was used for those samples in which the analyte was not detected at the detection limit.
[e] Not detected; value represents half the detection limit.
[f] Essential nutrient

Table 3-16
Summary of Radionuclide Activity in Surface Water, Salmon Site

Radionuclide	Frequency of Detection	Minimum Detectable Activity[a] (pCi/ℓ)[b]	Range of Detected Concentrations (pCi/ℓ)	Reference Range[c,d] (pCi/ℓ)	Average Concentration[d] (pCi/ℓ)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (pCi/ℓ)	Exceeds Reference Range?
Tin-113	1/7	7.10	8.70	3.90(ND)[e] - 5.50(ND)	6.02	1.86	7.74	No[f]
Tritium	5/12	0.19	0.20 - 0.70	0.10 - 0.20	0.22	0.19	0.34	Yes

[a]Minimum Detectable Activity (MDA) was selected from those samples in which the analyte was not detected.
[b]Picocurie per gram
[c]Reference range consists of samples GrC-1, HHC-1, HMC-1, and ReP-1.
[d]Half the MDA was used for those samples in which the analyte was not detected at the MDA.
[e]Not detected; values represent half the MDA.
[f]The only detection was lower than the actual MDA for several of the reference samples.

Table 3-17
Summary of Inorganic Concentrations in Surficial Soil, Salmon Site
(Page 1 of 2)

Analyte	Frequency of Detection	Minimum Detection Limit[a] (mg/kg)[b]	Range of Detections (mg/kg)	Reference Range[c,d] (mg/kg)	Average Concentration[d] (mg/kg)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (mg/kg)	Constituent of Potential Concern
Aluminum	108/108	--	719.00 - 18,800	1,980 - 14,600	4,694	2,883	5244.47	No[e]
Antimony	1/90	12.10	23.60	6.05(ND)[f] - 8.10(ND)	6.88	2.23	7.34	No[c,g]
Arsenic	106/108	0.54	0.49 - 11.90	0.66 - 2.70	1.88	1.44	2.15	No[e]
Barium	108/108	--	3.50 - 429	9.60 - 80.60	37.76	51.67	47.61	No[e]
Beryllium	18/108	0.06	0.07 - 0.29	0.03(ND) - 0.33	0.09	0.09	0.11	No[e]
Cadmium	7/108	0.64	0.72 - 1.40	0.32(ND) - 0.45(ND)	0.39	0.17	0.42	No[e]
Calcium	108/108	--	144 - 9,480	273 - 671	805.47	1,201	1,035	No[h]
Chromium	108/108	--	1.50 - 44.00	0.48(ND) - 10.40	7.74	5.90	8.89	No[e]
Cobalt	60/108	0.87	1.00 - 10.60	0.45(ND) - 5.60	2.58	2.11	2.98	No[e]
Copper	34/108	0.55	1.20 - 9,950	0.27(ND) - 10.70	97.20	957	280	Yes
Iron	108/108	--	1,170 - 19,700	1,780 - 11,200	6,390	3,721	7,100	No[e,h]
Lead	105/108	5.20	1.40 - 120.00	1.80 - 12.60	14.09	16.92	17.32	Yes
Magnesium	107/108	161	49.80 - 1,500	120 - 582	241	230	285	No[e,h]
Manganese	108/108	--	6.50 - 1,150	60.50 - 385	200	274	253	No[e]
Mercury	18/108	0.05	0.06 - 0.29	0.03(ND) - 0.04(ND)	0.04	0.04	0.05	No[i]
Nickel	10/108	4.30	4.70 - 19.10	2.11(ND) - 7.80	2.90	2.23	3.33	No[e]
Potassium	7/108	6.50	725 - 878	308(ND) - 656	393	307	453	No[e,h]
Selenium	6/104	0.21	0.24 - 0.69	0.11(ND) - 0.21(ND)	0.15	0.10	0.17	No[e]
Sodium	16/108	11.30	16.4 - 154	24.55(ND) - 57.43	22.68	23.37	27.13	No[e,h]

Refer to footnotes at end of table.

Table 3-17
Summary of Inorganic Concentrations in Surficial Soil, Salmon Site
(Page 2 of 2)

Analyte	Frequency of Detection	Minimum Detection Limit[a] (mg/kg)[b]	Range of Detections (mg/kg)	Reference Range[c,d] (mg/kg)	Average Concentration[d] (mg/kg)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (mg/kg)	Constituent of Potential Concern
Thallium	3/108	0.21	0.26 - 0.28	0.11(ND) - 0.14(ND)	0.12	0.03	0.13	No[g]
Vanadium	104/108	2.50	5.40 - 34.00	4.10 - 25.60	8.27	7.93	9.78	No[e]
Zinc	88/108	2.10	2.40 - 278	3.60(ND) - 24.40	14.14	27.78	19.44	No[e]
Cyanide	18/108	0.13	0.13 - 0.68	0.07(ND) - 0.27	0.11	0.10	0.13	No[e]

[a]Minimum detection limit from those samples with no detection
[b]Milligram per kilogram
[c]Reference range consists of samples Ref-19A, Ref-19B, Ref-19C, Ref-19D, and Ref-19E.
[d]Half the detection limit was used for those samples in which the analyte was not detected at the detection limit.
[e]95-percent upper confidence limit of the mean was within reference concentration range.
[f]Not detected; value represents half the detection limit.
[g]Detected in less than 5 percent of samples
[h]Essential nutrient
[i]95-percent upper confidence limit of the mean did not exceed the actual reference detection limits.

Table 3-18
Summary of Pesticide and PCB Concentrations in Surficial Soil, Salmon Site

Analyte	Frequency of Detection	Minimum Detection Limit[a] (µg/kg)[b]	Range of Detected Concentrations (µg/kg)	Reference Range[c,d] (µg/kg)	Average Concentration[d] (µg/kg)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (µg/kg)	Constituent of Potential Concern
4,4'-DDD	1/99	34	41.00	16.50(ND)[e] - 22.00(ND)	20.49	19.18	24.31	No[f]
4,4'-DDE	2/99	34	94.00 - 210	16.50(ND) - 22.00(ND)	22.97	27.91	28.53	No[f]
4,4'-DDT	2/99	34	90.00 - 190	16.50(ND) - 22.00(ND)	22.73	26.48	28.01	No[f]
Arochlor 1254	1/99	340	1,100	16.50(ND) - 220.00(ND)	212	211	254	No[f]
Dieldrin	3/99	35	42.00 - 320	16.50(ND) - 22.00(ND)	26.63	46.48	35.90	No[f]
Endosulfan II	1/99	35	38.00	16.50(ND) - 22.00(ND)	20.49	19.15	24.31	No[f]
Heptachlor epoxide	1/99	17	62.00	8.00(ND) - 11.00(ND)	10.88	11.12	13.09	No[f]

[a] Minimum detection limit from those samples with no detection
[b] Microgram per kilogram
[c] Reference range consists of samples Ref-19A, Ref-19B, Ref-19C, Ref-19D, and Ref-19E.
[d] Half the detection limit was used for those samples in which the analytes were not detected at the detection limit.
[e] Not detected; value represents half the detection limit.
[f] Detected in less than 5 percent of samples

Table 3-19
Summary of Semivolatile and Volatile Organic Concentrations in Surficial Soil, Salmon Site

Analyte	Frequency of Detection	Minimum Detection Limit[a] (µg/kg)[b]	Range of Detected Concentrations (µg/kg)	Reference Range[c,d] (µg/kg)	Average Concentration[d] (µg/kg)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (µg/kg)	Constituent of Potential Concern
Acenaphthylene	1/103	340	110	170(ND)[e] - 225(ND)	225.55	81.15	241.41	No[f]
Acetone	6/106	10.00	24.00 - 97.00	5.50(ND) - 8.50(ND)	21.42	18.32	24.95	No[g,h]
Benzoic Acid	51/103	1,600	39.00 - 870	99.00 - 570	690.14	578.05	803.14	No[g]
Bis(2-ethylhexyl)phthalate	55/103	340	38.00 - 1,700	170(ND) - 225(ND)	217.82	241.75	265.08	No[h]
Chloroform	6/96	2.00	4.00 - 7.00	2.50(ND) - 3.50(ND)	2.93	0.92	3.12	No[g]
Ethylbenzene	4/96	5.00	1.00 - 2.00	2.50(ND) - 3.50(ND)	2.75	0.36	2.82	No[f,g,i]
Methylene Chloride	9/105	5.00	1.00 - 10.00	2.50(ND) - 3.50(ND)	4.30	3.01	4.88	No[g,h]
Phenol	2/103	340	48.00 - 130	170(ND) - 225(ND)	229.06	83.23	245.33	No[f]
Styrene	1/97	5.00	25.00	2.50(ND) - 3.50(ND)	3.98	2.58	4.50	No[f,g,i]
Toluene	25/101	5.00	1.00 - 9.00	2.50(ND) - 3.50(ND)	2.67	0.88	2.84	No[h]
Xylene (total)	13/96	5.00	2.00 - 13.00	2.50(ND) - 3.50(ND)	4.01	1.85	4.39	No[g]

[a] Minimum detection limit was selected from those with no detections.
[b] Microgram per kilogram
[c] Reference range consists of samples Ref-19A, Ref-19B, Ref-19C, Ref-19D, and Ref-19E.
[d] Half the detection limit was used for those samples in which the analyte was not detected at the detection limit.
[e] Not detected; value represents half the detection limit.
[f] Detected in less than 5 percent of samples
[g] Log K_{ow} less than 3.5, unlikely to bioaccumulate (Garten and Trabalka, 1983)
[h] Common laboratory contaminant
[i] Within reference concentration range

Table 3-20
Summary of Radionuclide Activity in Surficial Soil, Salmon Site
(Page 1 of 2)

Radionuclide	Frequency of Detection	Minimum Detectable Activity[a] (pCi/g)[b]	Range of Detected Concentrations (pCi/g)	Reference Range[c,d] (pCi/g)	Average Concentration[d] (pCi/g)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (pCi/g)	Exceeds Reference Range?
Gross Alpha	74/97	5.86	5.52 - 37.90	4.24 - 12.20	10.60	5.95	11.80	No[e]
Gross Beta	97/97	--	6.78 - 31.20	9.65 - 15.60	14.24	4.04	15.05	No[e]
Bismuth-212	45/75	0.09	0.16 - 1.01	0.06 - 0.46	0.29	0.21	0.34	No[e]
Bismuth-214	1/1	--	0.05	NA[f]	NA	NA	NA	NA
Cesium-137	86/98	0.02	0.02 - 0.92	0.01(ND)[g] - 0.38	0.19	0.20	0.23	No[c]
Lead-210	46/66	0.28	0.04 - 2.04	0.20(ND) - 2.05	0.95	0.44	1.06	No[e]
Lead-212	63/86	0.02	0.20 - 1.86	0.01 - 0.53	0.47	0.41	0.56	Yes
Mercury-203	1/1	--	0.03	NA	NA	NA	NA	NA
Potassium-40	84/87	0.49	0.25-7.16	1.27 - 5.39	2.10	1.53	2.42	No[e]
Proactinium-231	5/5	--	0.61 - 0.96	NA	0.75	0.14	0.92	NA
Protactinium-234	1/1	--	2.96	NA	NA	NA	NA	NA
Radium-223	14/14	--	0.09 - 0.42	0.16	0.21	0.10	0.27	Yes
Radium-224	22/65	0.18	0.29 - 1.46	0.08 - 0.59	0.34	0.29	0.41	No[e]
Radium-226	79/87	0.03	0.19 - 1.29	0.41 - 0.73	0.47	0.27	0.53	No[e]
Radium-228	76/87	0.06	0.18 - 1.76	0.42 - 0.72	0.47	0.33	0.54	No[e]
Tin-113	2/2	--	0.02 - 0.03	NA	0.03	0.01	0.09	NA

Refer to footnotes at end of table.

Table 3-20
Summary of Radionuclide Activity in Surficial Soil, Salmon Site
(Page 2 of 2)

Radionuclide	Frequency of Detection	Minimum Detectable Activity[a] (pCi/g)[b]	Range of Detected Concentrations (pCi/g)	Reference Range[c,d] (pCi/g)	Average Concentration[d] (pCi/g)	Standard Deviation[d]	95% Upper Confidence Limit of the Mean[d] (pCi/g)	Exceeds Reference Range?
Thallium-208	53/80	0.01	0.07 - 0.53	0.11 - 0.15	0.13	0.12	0.16	Yes
Thorium-228	11/11	--	0.83 - 11.61	0.38(ND)	3.00	3.09	5.08	Yes
Uranium-234	2/2	--	3.19 - 10.40	NA	6.80	5.10	52.62	NA
Uranium-235	1/1	--	0.09	NA	NA	NA	NA	NA
Uranium-238	64/87	0.17	0.21 - 1.80	0.52 - 0.88	0.60	0.38	0.68	No[e]

[a]Minimum Detectable Activity (MDA) was selected from those samples in which the analyte was not detected.
[b]Picocurie per gram
[c]Reference range consists of samples GrC-1, HMC-1, HHC-1, and ReP-1.
[d]Half the MDA was used for those samples in which the analyte was not detected at the MDA.
[e]Maximum detected sample concentration was less than or equal to reference.
[f]Not applicable
[g]Not detected; value represents half the MDA.

Gamma spectroscopic analysis of sediment samples denoted several detectable levels of radionuclides. Although most radionuclides were detected infrequently and at reference concentrations, cesium-137 and lead-210 were found elevated in the Gator Pond and Beaver Pond sediments. In these ponds, potassium-40 was detected at 5.21 pCi/g and 2.38 pCi/g, cesium-137 at 0.61 pCi/g and 1.09 pCi/g, respectively, and lead-210 at 9.6 pCi/g in the Beaver Pond.

Tritium was detected in one sediment sample on the SS. This sample was from the Beaver Pond at a concentration of 0.48 pCi/g. Radioactive constituent data are summarized in Table 3-12.

Grain-size distribution data show the creek sediments to be at least 90-percent sand. Though the pond stations had smaller particle sizes, the majority was either sand or silt. Total organic carbon content ranged from less than 0.1 percent in Grantham Creek to over 19 percent in the Beaver Pond. Ammonia at the Beaver Pond station was extremely high (195 µg/kg); ammonia at other areas ranged from below detection to 3.43 µg/kg. These data are presented in Table 3-13.

3.2.3.2 Surface Water

Surface water data are presented in Appendix A. Tables A-4.1 through A-4.9 present the inorganic, pesticide and PCBs, volatile, semivolatile, gross alpha and beta, gamma spectroscopy, tritium, wet chemistry, and BOD data sets, respectively.

The majority of inorganic constituents detected in surface water are naturally occurring soil constituents. Many are essential macronutrients and necessary components of surface water bodies. Trace elements, such as cobalt, arsenic, and mercury, were detected once each in the three ponds on site: cobalt at 9.1 micrograms per liter (µg/ℓ) in the Half Moon Creek Overflow Pond, arsenic at 2.2 µg/ℓ in the Beaver Pond, and mercury at 0.17 µg/ℓ in the Gator Pond. A summary of the inorganic constituents in surface water on the SS are presented in Table 3-14.

The only radioactive constituents detected in surface water were tin-113 and tritium (Table 3-15).

3.2.3.3 Surficial Soil

Soil data are presented in Appendix A. Tables A-5.1 through A-5.7 present the inorganic, pesticide and PCBs, volatile, semivolatile, gross alpha and beta, and gamma spectroscopy data sets, respectively.

Many of the inorganic constituents detected are naturally occurring soil constituents and were found at reference levels. One copper concentration of almost 10,000 mg/kg was found in surficial soils at the West Substation although most other samples were comparable to reference (ND - 10.70 mg/kg) and background concentrations (ND - 700 mg/kg). Concentrations of inorganics in surficial soils on the SS are summarized in Table 3-17.

Several pesticides were detected in soils on site at isolated locations. The Big Chief Drilling Storage Area in SA-3 contained detectable concentrations of heptachlor epoxide (62 µg/kg), dieldrin (42 and 320 µg/kg), and endosulfan II (38 µg/kg). Dieldrin was also detected at the North Substation in SA-4 (320 µg/kg). Both 4,4'-DDE and 4,4'-DDD were found at the East Substation in SA-1 (210 and 41 µg/kg, respectively) and the HT-2 Well Area in SA-5 (94 and 90 µg/kg, respectively). In addition, 4,4'-DDT was found at the East Substation (190 µg/kg).

One detection of PCBs was found in surficial soils. The East Substation had a detectable level of Aroclor-1254 (1,100 µg/kg). Concentrations of pesticides and PCBs in surficial soils in the SS are summarized in Table 3-18.

Volatile organic compounds detected in surficial soils included toluene, xylene, and styrene. Detected concentrations of toluene ranged from 1 µg/kg to 9 µg/kg, and detected concentrations of xylene ranged from 2 µg/kg to 13 mg/kg. The majority of detected concentrations were found at Station 1A, Big Chief Drilling Storage Area, Sites E-2 and E-7 in SA-3, and the CH Fuel Storage Area in SA-4.

Four semivolatile organic compounds were detected in surficial soils. Benzoic acid and bis(2-ethylhexyl)phthalate were detected in approximately half of the samples. Concentrations ranged from nondetectable to 1,900 µg/kg for benzoic acid and from nondetectable to 1,700 µg/kg for bis(2-ethylhexyl)phthalate. Phenol and acenaphthylene were detected twice and once, respectively. Concentrations of volatile and semivolatile organics in surficial soils in the SS are summarized in Table 3-19.

Gross alpha and beta and gamma spectroscopic data for surficial soils are summarized in Table 3-20. Although most sample concentrations were within reference ranges, elevated concentrations existed on site. Maximum on-site concentrations were 37.9 pCi/g and 31.2 pCi/g, respectively.

3.2.4 Constituents of Potential Concern

Constituents of potential concern are site related, have reported quantities of reliable quality, and are of greatest potential concern for ecological receptors given their physiochemical, toxicological, and bioaccumalative properties. A summary of the screening criteria is presented in Figure 3-3. It is this list of COPC for which stress was characterized at the SS.

3.2.4.1 Sediment

In sediment, six inorganic constituents, arsenic, barium, cadmium, lead, manganese, and zinc, were considered to be COPC. No organic constituents were considered COPC in sediment. Alpha and beta emissions are considered COPC in all media. Likewise, all radionuclides detected with sufficient quantity and quality are considered potentially site related and COPC. Tritium is considered a COPC because of past site activities. Three radionuclides were detected by gamma spectroscopy at levels above reference. These are cesium-137, lead-210, and radium-224.

3.2.4.2 Surface Water

In surface water, two inorganics are considered COPC: aluminum and arsenic. Aluminum was not considered a COPC in other media; however, arsenic was considered a COPC in sediment. No organic constituents were considered COPC in surface water.

Alpha and beta emissions were not detected in surface water samples. Tritium was detected in five samples. The highest levels were in the Beaver Pond and Half Moon Creek Overflow Pond. The only radionuclide detected by gamma spectroscopy in surface water samples was tin-113 in Grantham Creek just above the detection limit.

3.2.4.3 Surficial Soil

In soil, only two inorganic constituents are considered COPC. These are copper and lead. The majority of inorganic constituents detected were within site-specific reference ranges. Antimony, selenium, and thallium were detected in less than 5 percent of the surficial soil samples. None of these is attributable to past site activities.

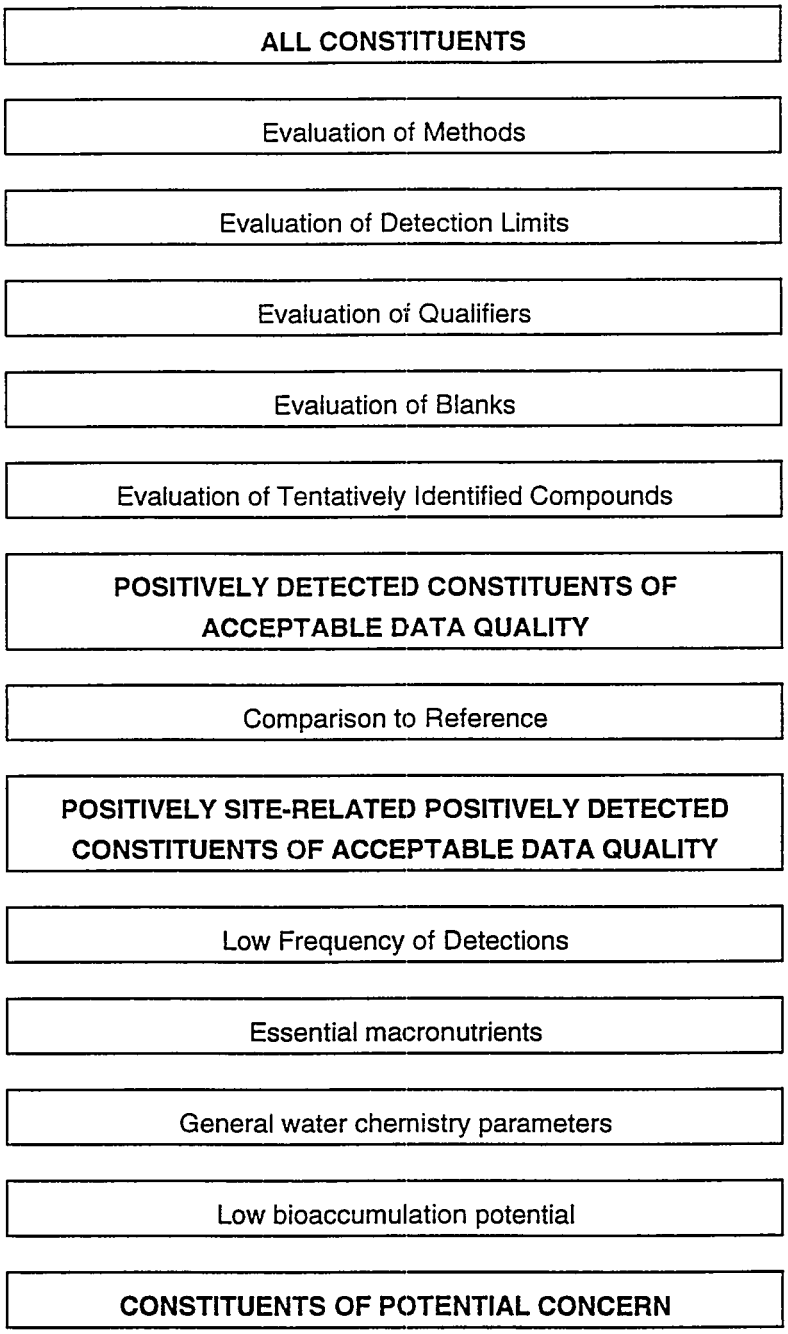

Figure 3-3
Screening Criteria for Selection of Constituents of Potential Concern

No organic constituents are considered COPC in soil. Phenol, acenaphthylene, ethylbenzene, and styrene were detected in less than 5 percent of the samples. Bis(2-ethylhexyl)phthalate is a common laboratory contaminant with no site-related history. Benzoic acid is a naturally occurring constituent found in both flora and vertebrate excretions and, along with toluene and xylene, have log K_{ow} values of less than 3.5.

Alpha and beta emissions are considered COPC in all media. Likewise, all radionuclides detected with sufficient quantity and quality through gamma spectroscopy are considered potentially site-related and COPC.

3.3 Exposure Pathway Identification

This pathway assessment analyzes potential exposure pathways for terrestrial and aquatic flora and fauna at the SS, using existing site conditions, coupled with additional information on the life histories of biota using the site. A Conceptual Site Model (CSM) (Figure 3-4) was used to describe suspected sources and types of constituents present, constituent release and transport mechanisms, rate of constituent release and transport (where possible), affected media, known and potential routes of migration, and known and potential environmental receptors.

Complete exposure pathways must exist for exposures to occur. A complete exposure pathway requires (1) a source and mechanism for contaminant release, (2) a transport medium, (3) a point of environmental contact, and (4) an exposure route at the exposure point (EPA, 1989b). If any of these four components is absent, a pathway is generally considered incomplete. However, the transport medium may be missing, and the pathway may still be complete if the contact point is directly at the contaminant release point.

Contaminants at the SS may migrate from their source to other environmental compartments via six major pathways: (1) volatilization to the air and transport by wind, (2) leaching to the saturated zone and transport with groundwater, (3) adsorption onto surface soil particles and transport by wind as fugitive dust, (4) erosion and transport with surface water runoff, (5) discharge of groundwater to surface water, and (6) exchanges between surface water and sediment.

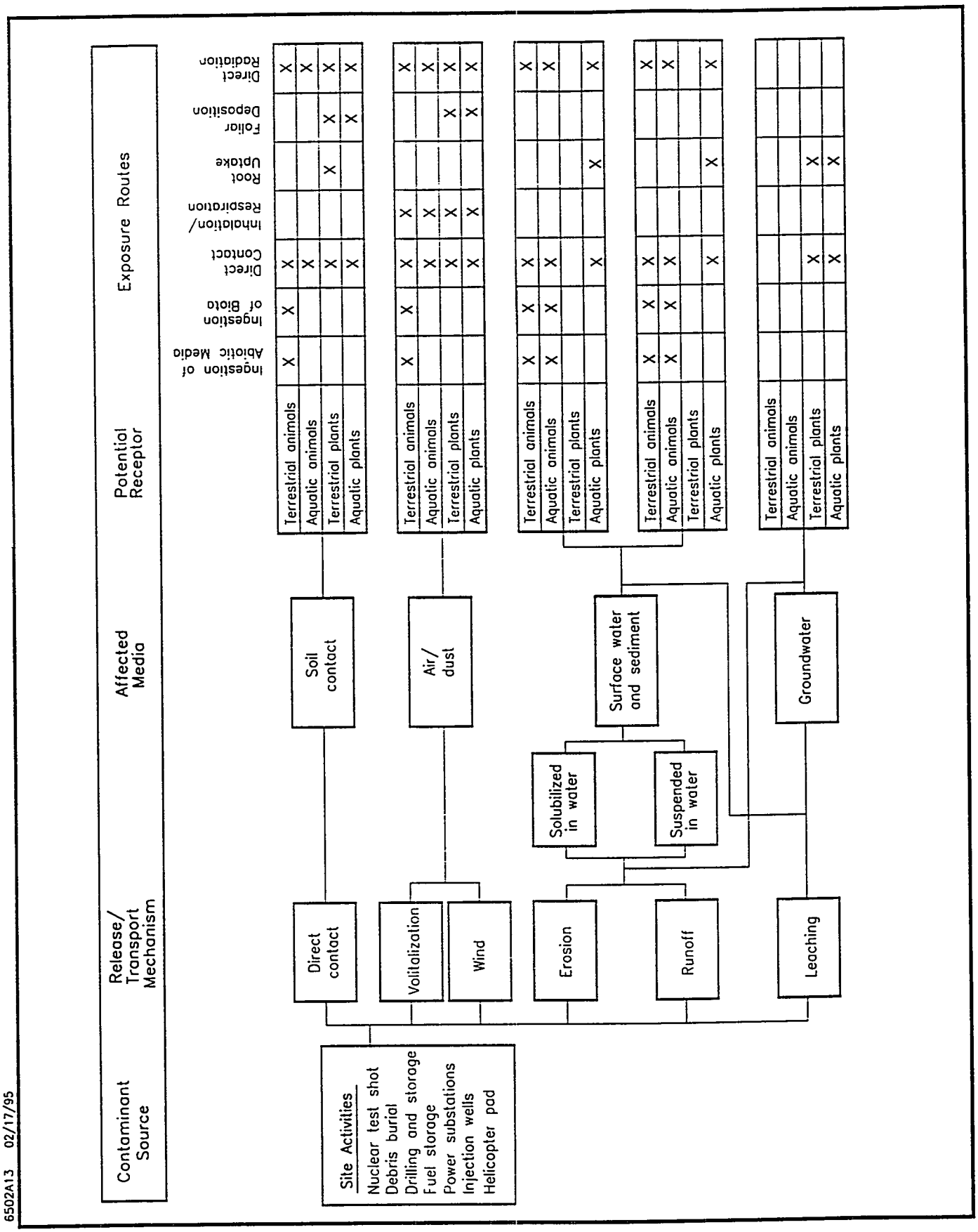

Figure 3-4
Conceptual Site Model,
Salmon Site

Nonadsorptive water-soluble contaminants present in soils may migrate into groundwater or become dissolved in stormwater runoff from the site. Contaminants that leach into the groundwater may be transported to surface waters via seeps and/or springs. Contaminants with strong affinity for soils may be transported off site as sediment load in surface water or as wind-generated particulates. Nonvolatile contaminants with a lesser affinity for soils may also be transported off site by wind.

Ecological receptors may be exposed to water-borne contaminants via dermal contact, consumption of water, or inhalation of organic vapors. Terrestrial organisms may come into contact with water-borne contaminants as a result of wading or swimming in contaminated waters. Terrestrial organisms may also ingest water-borne contaminants if wildlife use impacted waters as a source of drinking water. Aquatic organisms inhabiting contaminated waters would be in constant contact with contaminants. Terrestrial organisms may be exposed to organic vapors that result from the volatilization of organics in surface waters; however, significant exposure via this route would be limited to terrestrial organisms that spend most of their time near or in the water (e.g., waterfowl and turtles).

Contaminants present in sediment may result from erosion or adsorption of water-borne contaminants onto sediment particles. If sediments are present in an area that is periodically inundated with water, previous exposure pathways for soils would be applicable during dry periods. Water overlying sediments would prevent contaminants from either volatilizing or being carried by wind erosion. Exposure via dermal contact may occur especially for benthic organisms. Some aquatic organisms consume sediment and ingest organic matter from the sediment. Inadvertent ingestion of sediments may occur as the result of feeding on benthic organisms and plants.

While contaminants in soils may leach into groundwater, environmental receptors generally will not come in contact with these contaminants unless groundwater is discharged to the surface. However, root contact with, and uptake of, contaminants in groundwater is a potential pathway.

Potential primary exposure pathways for terrestrial receptors include (1) dermal contact with contaminated soil, sediment, and surface water; (2) ingestion of contaminated soil, sediment, and surface water; (3) inhalation of organic vapors and/or contaminated particulates; (4) uptake through roots in contact with soil, sediment, surface water, or groundwater; and (5) direct radiation from contaminated soil, sediment, or surface water.

Potential primary exposure pathways for aquatic receptors include (1) dermal contact with contaminated sediments or surface water, (2) ingestion of contaminated sediment or surface water, (3) uptake through roots and leaves in contact with surface water or sediment, and (4) direct radiation from contaminated surface water or sediment.

Secondary exposure pathways involve contaminants that are transferred through the food chain. This pathway may include contaminants bioaccumulated from soil or sediment into plant tissues, which are subsequently ingested by terrestrial or aquatic herbivores. Benthic invertebrates may bioaccumulate contaminants from the surface water and/or sediments and then become prey for bottom feeding fish, which subsequently become prey for piscivorous avian species.

The potential sources of contamination and routes of migration are summarized in the site-wide CSM for the SS (Figure 3-4).

3.3.1 Animal Exposure Routes

There are five major exposures routes by which terrestrial and aquatic animals might be exposed to contaminants at the SS: (1) direct ingestion of contaminated abiotic media, (2) the consumption of contaminated animal or plant tissue (food web interactions), (3) inhalation, (4) direct contact (e.g., absorption through skin or gill surfaces), and (5) direct radiation.

Direct Ingestion of Contaminated Abiotic Media

Direct ingestion of contaminated soil or sediment could occur while animals grub for food, feed on plant matter covered with contaminated soil, filter feed in areas where sediments have been resuspended in the water column, or preen or groom themselves. Soil ingestion may inadvertently occur while grooming; burrowing; or consuming plants, insects, or invertebrates resident in the soil. Aquatic deposit feeders directly ingest large quantities of bulk sediment in order to obtain the energy-rich organic fraction; these organisms would likely have a significant exposure from this pathway. Terrestrial species would also ingest potentially contaminated surface water from the streams, ponds, and wetland areas at the SS.

Ingestion of Contaminated Tissues

Predatory organisms (secondary and tertiary consumers) may be exposed when feeding upon prey that contains elevated levels of contaminants. The ability of contaminants to be passed from lower to higher trophic levels is dependent upon their fate inside the prey animals. If the contaminant is solely ingested and excreted by the prey without absorption from the gastrointestinal (GI) tract and storage in its tissues, the exposure of higher trophic levels is

minimized. The dose received by the predator is dependent upon the contaminant concentration in the prey, its ingestion rate, and the amount of contaminated matter consumed.

The risk of exposure for predatory organisms is greater if the prey assimilates the ingested contaminants or directly bioaccumulates contaminants present in the media into its tissues. The risk of exposure to predators in this case would then be dependent upon the concentration of contaminants in the particular tissues consumed and the rate of food consumption. The actual concentration within the predator would also depend upon the rate of assimilation from the GI tract of the contaminants during ingestion, the rate of metabolism, and the depuration rate.

Inhalation

Volatile organic compounds and, to a lesser extent, lower-molecular-weight polycyclic aromatic hydrocarbons tend to volatilize from surface soils or surface water. In vapor form, these compounds may become bioavailable to organisms during respiration and pose an important exposure route. However, organic compounds were not detected in high concentrations in environmental media, and this exposure route is probably minimal at the SS.

Exposure via inhalation of fugitive dust would be limited to contaminants present in surface soils at an area that was devoid of either vegetation or landscaping that would prevent the erosion of soil particles. Exposure via inhalation is most important to organisms that burrow in contaminated soils, especially those that bear their young in below-ground burrows (e.g., gopher tortoises). However, intake via this mechanism is considered minimal relative to other pathways.

Direct Contact

Direct exposure to soils, sediments, and surface water are other pathways important to some species. This could result from direct dermal contact with the soils, sediments, or surface water on unprotected surfaces (e.g., gill membranes, exposed skin, or exposed mucosal membranes). Significant exposure via dermal contact would be limited to organic contaminants that are lipophilic and can penetrate epidermal barriers. Mammals and birds are less susceptible to exposure via dermal contact with soils because their fur and feathers prevent skin from coming into full contact with soil.

Dermal exposure may be most important to aquatic organisms, such as benthic invertebrates, and bottom dwelling fish. Direct exposure is likely to have far less impact on terrestrial organisms than other routes such as ingestion of contaminated media or tissues.

Direct Radiation

Unlike chemicals, radionuclides can have deleterious effects on animals without being taken or brought into contact with the body. This is because high-energy beta particles from radionuclides in contaminated air, water, or soil can travel long distances with only minimum attenuation in these media before depositing their energy in animal tissues. Gamma and x-rays are the most penetrating of the emitted radiations and comprise the primary contribution to the radiation dose from external exposures. Alpha particles are not sufficiently energetic to penetrate the outer layer of skin and do not contribute significantly to the external dose. External exposure to beta particles primarily imparts a dose to the outer layer skin cells, although high-energy beta radiation can penetrate into animal bodies.

3.3.2 Plant Exposure Routes

There are four major routes by which plants might be exposed to contaminants at the SS: (1) incorporation of organic vapors through gas exchange, (2) root uptake, (3) foliar deposition or direct contact for aquatic plants, and (4) direct radiation.

Respiration of Organic Vapors

Organic vapors resulting from the volatilization of contaminants in soil and/or surface water may be incorporated into vascular plants via gas exchange and respiration. Because the SS did not have high media levels of organics, exposure via this pathway is probably minimal.

Root Uptake

Both aquatic and terrestrial plants may be exposed to site-related contaminants in soil, sediment, and/or surface water by root uptake. This pathway is the most important exposure route for plant uptake of contaminants.

Plant uptake of chemicals and radionuclides from soils is affected by numerous processes. Because of the multiplicity of factors, soil-to-plant concentration ratios may exhibit considerable variability. Among the more important factors that affect root uptake are (1) the physicochemical form of the chemical or radionuclide, (2) soil characteristics, (3) fertilizers and agricultural chemicals, (4) chelating agents in the soil, (5) distribution of chemicals and radionuclides in the soil, and (6) plant species and internal translocation within the plant.

Foliar Deposition

Plants on and around the SS may be exposed to site-related contaminants adhering to particulates that have been eroded by the wind. Wet and/or dry deposition of air-associated contaminants

onto leaf surfaces would make them available for uptake by the plants either directly through the epidermal tissues or through the stomates. Uptake of contaminants may also occur through direct contact of aquatic plants with contaminated surface waters.

Direct Radiation

High-energy beta particles from radionuclides in contaminated air, water, or soil can travel long distances with only minimal attenuation in these media before depositing their energy in plant tissues. Gamma and x-rays are the most penetrating of the emitted radiations and comprise the primary contribution to the radiation dose from external exposures. Alpha particles are not sufficiently energetic to penetrate the outer layer of plant tissue and do not contribute significantly to the external dose. External exposure to beta particles primarily imparts a dose to the outer layer of plant tissue although high-energy beta radiation can penetrate further into plant tissues.

The buildup of radionuclides on vegetation as a result of deposition of airborne radioactive materials may be an important process in several exposure pathways. It is a major source of radionuclide contamination of terrestrial food products such as vegetables and forage for terrestrial animals. Deposited beta- and gamma-emitting radionuclides contribute to exposure from external radiation.

3.4 Ecological Receptor Identification

An ecosystem is the combination of the abiotic (nonliving) physiochemical environment and the assemblage of biotic (living) organisms that combine together to form an interrelated and interdependent system.

There are two basic environments existing at SS that provide a natural classification of chemical and/or radiation exposure pathways: terrestrial and aquatic (freshwater) ecosystems. For the SS ecosystems, the areal extent of investigation has been defined as the former test site bounded by the fence. However, it should be noted that the habitats present on site (e.g., forest, field, and wetlands) extend off-site beyond the area of investigation.

The loss of valued native species or the establishment of unwanted exotic species would be viewed as a major ecosystem degradation. Closely tied to this view is the concept and measure of species' diversity. Diversity has been shown to generally decline in response to disturbance for a variety of different stressors and ecosystems.

3.4.1 Terrestrial Ecosystems

3.4.1.1 Vegetation

The SS site is dominated by open long-leaf pine forest (*Pinus taeda* and *Pinus palustris*) with a shrub understory. Some areas have been replaced by oaks (*Quercus* sp.), hickory (*Carya* sp.), and magnolia (*Magnolia* sp.).

Along the drainages and in the vicinity of the ponds on the site, bottomland hardwood forest with bald cypress (*Taxodium distichum*), water tupelo (*Nyssa aquatica*), red maple (*Acer rubrum*), black willow (*Salix nigra*), and other obligate and facultative wetland tree species comprise the canopy and understory. Emergent wetland vegetation exists around open ponded areas and is composed of water lily (*Nymphaea odorata*), common cattail (*Typha latifolia*), buttonbush (*Cephalanthus occidentalis*), marsh elder (*Iva frutescens*), and arrow arum (*Peltandra virginica*) among other species.

3.4.1.2 Animals

Primary consumers at the site include large mammals (white-tailed deer [*Odocoileus virginianus]*) and small mammals (mice [*Peromyscus* sp.], short-tailed shrews [Blarina brevicauda], pine vole [*Pitymus pinetorum*], and eastern cottontail [*Sylvilagus floridanus]*). Secondary consumers include omnivorous raccoons (*Procyon lotor*), opossums (*Didelphis marsupialis*, and carnivorous coyotes[*Canis latrans*]).

Birds identified at SS are mostly passerines including the Bachman's sparrow (*Aimophila aestivalis*), which is on the threatened species list, and the northern bobwhite (*Colinus virginianus*), which is a common game bird. A number of avian predators were also identified, including the red-tailed hawk (*Buteo jamaicensis*).

Reptiles and amphibians can be found in both aquatic and terrestrial habitats. Most of these animals at the site were observed in aquatic environments. However, a terrestrial sensitive species, the gopher tortoise (*Gopherus polyphemus*), was observed at the site.

3.4.2 Aquatic Ecosystems

3.4.2.1 Macrobenthos

The qualitative macrobenthos field data are summarized in Table 3-21. Ten taxa were observed at the various creek stations. Taxa identification and confirmation were performed using

Peckarsky et al., 1990. The most prevalent taxa found throughout the site were Plecoptera, Coleoptera, Anisoptera, Decapoda, and Tabanidae. Station HHC-2 was the lowest in overall taxa richness, but was abundant in Chironomidae, which are generally a stress- tolerant species. This was probably more attributable to the station's creek topography than external sources of stress. Station HMC-3 had a large submerged root mass present that was found to support a diverse macroinvertebrate population. Grass shrimp and gastropods, taxa usually indicative of a healthy macrobenthos population, were unique to that station, which is located in close proximity to SGZ.

Stations HMC-2, HMC-5, and GRC-2 shared a common overall abundance and taxa richness, with three different taxa present. Station HHC-2 had two taxa that were abundant (>10 individual organisms) and three taxa that were rare (<3 individuals). Stations HHC-1, GRC-1, HMC-1, HMC-3, and HMC-4 shared a common overall abundance and taxa richness, with four to seven different taxa occurring at those stations. Crayfish, lampreys, salamanders, tadpoles, Anisoptera, Coleoptera, and Plecoptera were found commonly throughout the creek areas. The presence of these taxa indicates that a relatively high-quality habitat is present throughout the creek systems, with no observable site-related influence.

3.4.2.2 Reptiles and Amphibians

Salamanders are commonly found in or near creeks, springs, or seeps and were found at Half Moon Creek on the site. They are usually absent from large streams where predatory fish live. They are most common along the edges of small woodland streams where stones, chunks of wood, and miscellaneous debris provide ample shelter and food. Salamanders feed on a number of aquatic insects and other aquatic macroinvertebrates. They are in turn fed upon by larger salamanders, fish, birds, and mammals. Aquatic amphibians are good indicators of the health of aquatic systems. These animals are especially sensitive to pollution and loss of aquatic habitat. Snakes, including a water snake (*Nerodea* sp.), were observed along with lizards; turtles; skinks; toads; and frogs, including the river frog (*Rana heckscheri*).

3.4.2.3 Finfish

Fifteen finfish species were found on the site. The greatest diversity and abundance occurred at station HMC-4. No fish species were observed or collected at stations HHC-1 or GtP-1. The only abnormalities observed were external parasites on a bluegill (*Lepomis macrochirus*) found at station GRC-1.

Table 3-21
Macrobenthos Qualitative Field Data,
Salmon Site

Taxon	Salmon Site Station Locations								
	HMC-1	HMC-2	HMC-3	HMC-4	HMC-5	GRC-1	GRC-2	HHC-1	HHC-2
Oligochaeta	[a]								
Isopoda									
Amphipoda									
Decapodaa									
Insecta	C[b]	C	C	R[c]				C	R
Anisoptera	C	C	C	R	C	C	C	C	A[d]
Zygoptera	C		C						
Hemiptera									
Coleoptera	C		C		C	A	A	A	R
Diptera									
Tipulidae									
Tabanidae	R			C		R		C	
Culicidae						C			R
Chironomidae	C		C	A	C	C	C		
Plecoptera	C	C							
Ephemeroptera									
Tricoptera									
Gastropoda			C						

[a] Blank = Not Observed
[b] Common 3-9 Individuals
[c] Rare < 3 Individuals
[d] Abundant > 10 Individuals

The cherryfin shiner (*Lythrurus roseipinnis*) was found at all three creeks and was the dominant species on the site. It was dominant at creek stations GRC-1, GRC-2, HMC-2, HMC-3, HMC-4, and HMC-5 and was also found at creek stations GRC-1, HMC-2, and HMC-3. The two other shiner species found on the site were the silverstripe shiner (*Notropis stilbuis*) at pond station BeP-1 and creek station HMC-4 and the blacktail shiner (*Notropis venusta*) found only at creek station HMC-4. Two darter species were found on the site: the blackbanded darter (*Percina nigrofasciata*) at creek station GRC-1 and the gulf darter (*Etheostoma swaini*) at the creek station HMC-3.

Additional fish species found on the site were the starhead top minnow (*Fundulus notti*), which was found at pond stations BeP-1 and BeP-2 and creek station HMC-4. This species was dominant at BeP-2. The chestnut lamprey (*Ichthyomyzon castaneus*) was found at three creek stations: GRC-2, HMC-4, and HMC-5. The bluegill (*Lepomis macrochirus*) was dominant at pond station HOP-1 and was also found at creek station GRC-1. The mosquitofish (*Gambusia affinis*) was dominant at pond station BeP-1 and was also found at pond station BeP-2.

3.5 Potential Ecological Effects

Chemical and radiological contaminants are the primary stressors at the SS. Other anthropogenic, physical, or naturally-occurring stressors are discussed in this assessment where appropriate; however, potential impacts to ecological receptors at the SS from significant off-site (non-SS) stressor sources were not investigated for this BERA.

For a given set of environmental conditions, species have characteristic attributes, such as birth rates, age and sex distributions, migration patterns, and mortality rates. A species' habitat preferences, food preferences, and other behavioral characteristics (e.g., nesting, foraging, and rearing young) also may determine population size and distribution in an area and may also significantly affect the potential for exposure.

A contaminant entering the environment will cause adverse effects if

- It exists in a form and concentration sufficient to cause harm
- It comes in contact with organisms or environmental media with which it can interact
- The interaction that takes place is detrimental to life functions.

Adverse effects may also occur if a contaminant interacts with other chemicals already present in a synergistic manner that could raise the overall toxicity of the contaminated environment. The likelihood of ecological harm is thus a combined function of chemical, physical, and biological

factors, depending on the nature of the contaminant and the environment into which it is released.

Ecological effects from a single or occasional release are likely to be considerably different than those associated with a continuous release. Contaminants may enter the environment or move among environmental compartments on several possible time scales, such as the following:

- Single event (e.g., from an accidental spill)
- Intermittently (e.g., from storms transporting surficial soil runoff)
- Seasonally (e.g., from heavy rainfall during the spring)
- Continuously (e.g., from groundwater discharge).

Some of these release scenarios may be occurring at the SS, and each type of release may produce a different level of stress on a given ecosystem. Since the AEC's testing program at the SS ceased in 1970, the only contaminant releases to the SS environment potentially occurring are chronic in nature (e.g., groundwater discharge) or the result of intermittent episodes (e.g., heavy rainfall events). The assessment endpoints and measurement endpoints that have been selected for this assessment will address these different time spans and exposure durations, as well as the synergistic or antagonistic effects of different combinations of contaminants.

As stated above, the potential exposures to SS-related contaminants are expected to be chronic in nature. Chronic toxicity due to exogenous chemicals in an ecosystem can greatly increase the mortality rate of component populations or can change the organisms' ability to survive and reproduce in less direct ways, such as the following:

- Altering developmental rates, metabolic processes, physiologic function, or behavior patterns
- Increasing susceptibility to disease, parasitism, or predation
- Disrupting reproductive functions
- Causing mutations or otherwise reducing the viability of offspring.

Ecological effects are most easily expressed as some impairment of a biological function or condition. There are two basic approaches to expressing ecological effects: (1) observed effects or a statement of contaminant-related effects observed on site and an interpretation of associated ecological implications in relation to appropriate endpoints or (2) comparison of on-site measured concentrations to established benchmarks for the environmental media of concern. Both of these approaches, observed effects and comparison to benchmarks, were used in the ecological assessment of the various ecosystems at the SS.

Contaminants may cause or induce changes in the composition and structure of a biotic community as a secondary effect of the changes in the size of particular populations. These species may be a major source of food or shelter for the rest of the community. Others may be crucial in maintaining a balance of species in a habitat.

3.6 Endpoints

An assessment endpoint is an expression of the environmental value that is to be protected while the measurement endpoint is a measurable ecological characteristic that is related to the characteristic chosen as the assessment endpoint (Suter, 1993). The quality of the terrestrial and aquatic ecosystems at the SS and their vitality as habitat is the primary assessment endpoint for this study.

3.6.1 Aquatic Ecosystems

The aquatic ecosystems at the SS were assessed using two assessment endpoints as presented in Table 3-22: (1) Benthic macroinvertebrate community viability and (2) Probability of subacute affects in pelagic organisms. The quality and suitability as habitat of the streams and ponds that make up the aquatic ecosystems at the SS, namely Half Moon Creek, Grantham Creek, Hickory Hollow Creek, Beaver Pond 1-A, Beaver Pond 2, Gator Pond, and Half Moon Creek Overflow Pond, can be measured through an analysis of the benthic invertebrate community and the application of specific community metrics, as employed by RBP III (EPA, 1989c). Biological communities reflect overall ecological integrity, integrating chemical, physical, and biological influences, as well as the effects of different ecological stressors. Communities provide a holistic measure of the aggregate impact of these stressors over time and fluctuating environmental conditions.

Benthic invertebrates comprise a heterogenous assemblage of animal taxa that inhabit the sediment or live on or in other bottom substrates in the aquatic environment and are common in most aquatic habitats. The macroinvertebrate communities are good indicators of localized conditions because many invertebrates have limited migration patterns or sessile modes of life and are, therefore, well suited for assessing site-specific impacts. The macroinvertebrate communities also integrate the effects of short-term environmental variations.

Table 3-22
Assessment and Measurement Endpoints for Aquatic Ecosystems, Salmon Site

Assessment Endpoint	Measurement Endpoint
Benthic Macroinvertebrate Community Viability	(1) Rapid Bioassessment Protocol III Metrics. (2) Calculation of Estimated Aquatic Exposure Point Concentrations for Contaminants and Comparison to Literature-Derived Sediment or Water Quality Criteria.
Probability of Chronic Effects in Pelagic Organisms	(1) Fish Species Survey and Determination of Tissue Contaminant Burdens Relative to Literature-Derived Reference Levels. (2) Application of Ambient Water Quality Criteria. (3) Surface Water Toxicity Tests.

The measured community metrics from the streams and ponds associated with contaminant source areas at the SS are compared with the benthic macroinvertebrate community metrics of the reference locations to determine if site-related contaminants are impacting the ponds and streams at the SS and to determine the current overall biotic quality of the aquatic ecosystems.

In order to characterize the risk to the benthic macroinvertebrate communities, the estimated interstitial water concentrations of the COC were compared to ambient water quality criteria or standards for the protection of aquatic life or, when criteria were not available, threshold values as determined from the scientific literature.

The toxicity and kinetics of inorganics in sediment are not as well understood as for organics. For this reason, characterizing the risk to benthic organisms from metals in sediments will be accomplished by comparing the measured sediment concentration of the metals of concern to threshold concentrations as determined in the literature. The primary source of the biologically effective sediment concentrations is Long and Morgan (1990). Long and Morgan have compiled data from existing studies that link concentrations of contaminants with predicted or observed biological effects. From these data, concentrations were selected that represent the lower tenth percentile of the data. This concentration is the Effects Range - Low and is used as the benchmark value in the risk characterization for metals in sediments.

The second assessment endpoint was the probability of chronic effects in pelagic organisms in the ponds and streams at the SS. Chronic exposure to site COPC may induce deoxyribonucleic acid (DNA) damage, cancer, reproductive failure, or other effects. Fish populations were surveyed and characterized in the ponds and streams at the SS. Tissues were analyzed from several species of fish collected at the site. These body burdens can be compared with literature-derived effect levels to address the probability effects associated with chronic exposure. Surface water concentrations of COC were also measured in the ponds and streams at the SS, and comparisons were made with ambient water-quality criteria or standards for the protection of aquatic life. When these criteria were unavailable, comparisons were made with threshold values as determined from the scientific literature.

Comparisons were also made between fish tissue samples collected at reference locations that were not directly impacted by the SS and fish tissue samples collected from areas of suspected or known contamination.

3.6.2 Terrestrial Ecosystems

The quality and suitability as habitat of the terrestrial ecosystems at the SS were assessed using two assessment endpoints as presented in Table 3-23: (1) change in floral community diversity and (2) probability of chronic effects in faunal community.

The maintenance of diversity of a rich floral community is the first assessment endpoint to be considered for terrestrial systems at the SS. The floral communities throughout the SS were surveyed in order to identify individual species and communities present and their relative abundance. Visual evidence of stress or vitality was also noted during this survey. Information collected during these field surveys was used to characterize the vegetative communities and to determine whether they exhibited characteristics of either stressed or healthy communities.

Sampling and analysis of plant tissues from reference sites and source areas were also conducted at the SS. The analytical results from this sampling yields information regarding the potential plant uptake of site-related contaminants and their potential adverse effects on individual plants and communities. The concentrations of contaminants detected in samples from reference areas were compared to samples collected in suspected source areas to determine whether there were any significant differences in plant tissue concentrations. Literature-derived toxicity data are used to determine whether the increased tissue concentrations could cause detrimental effects to individual plants and/or communities.

The second assessment endpoint to be considered is the probability of chronic effects in the faunal communities at the SS. Chronic exposure to site COPC may induce DNA damage, cancer, reproductive failure, or other effects. The measurement endpoints will yield species-specific data that can be related to the entire SS. Avian species were surveyed in order to enumerate the different species using the site for nesting, breeding, feeding, or other activities.

Table 3-23
Assessment and Measurement Endpoints for Terrestrial Ecosystems, Salmon Site

Assessment Endpoint	Measurement Endpoint
Change in Flora Community Diversity	(1) Qualitative Evidence of Vegetative Stress or Vitality. (2) Vegetation Species Survey and Determination of Tissue Contaminant Burdens Relative to Literature-Derived Effects Levels.
Probability of Chronic Effects in Fauna Community	(1) Small Mammal Species Survey and Determination of Tissue Contaminant Burdens Relative to Literature-Derived Reference Levels. (2) Large Mammal Species Survey and Determination of Tissue Contaminant Burdens Relative to Literature-Derived Reference Levels. (3) Calculation of Estimated Exposure to Species Not Sampled (including Predators and Sensitive Species) and Comparison with Literature-Derived Effects Levels. (4) Terrestrial Toxicity Tests.

Small and large mammal surveys and tissue sampling were also used to aid in estimating the probability of adverse effects in these populations. These mammal surveys identified and enumerated the specific mammalian species utilizing the SS for feeding, breeding, hunting, and other activities. The analytical results of the mammalian tissues collected from the SS were used to identify whether site-related contaminants are bioaccumulating in mammals that frequent the site, and if so, in what specific tissue has the bioaccumulation taken place and to what extent.

3.7 Problem Formulation Summary

The SS can be segregated by ecosystem type into terrestrial and aquatic. Both ecosystems have COPC, receptors, and complete pathways. There is abundant habitat, and diverse plant and animal communities were observed in both ecosystems, but the potential exists for acute or chronic toxic effects from exposure to contaminants in soils, surface water, and sediments.

There are larger questions relative to the assessment goals for the SS: Are the COPC from previous site activities present in sufficient bioavailable concentrations, and are the COPC capable of bioaccumulation in SS flora and fauna to a level that might pose a risk to these organisms and these ecosystems? These questions were addressed via the analysis of the assessment and measurement endpoints through the processes of exposure effects and risk characterization in the following sections.

4.0 Exposure Characterization

Exposure characterization involves the quantification of the magnitude or type of actual and/or potential exposures of ecological receptors to site-specific COPC. This part of the assessment includes quantification of stressor release, transport and fate, ecological receptor characterization, and determination (either by measurement or modeling) of EPCs.

4.1 Transport and Fate Estimation

The EPC in sediment, surface water, and surficial soil samples are represented by the 95-percent UCL of measured concentrations of constituents that passed the screening process outlined in Figure 3-3, calculated as shown in Section 3.2.2.9. Additionally, soil/water partition coefficients (K_d), when available for inorganic COPC, were used to calculate aquatic EPC (EPC_{aq}) to estimate the concentration of COPC in interstitial waters as that portion of the bulk sediment concentration that was bioavailable. This was calculated by the following equation:

$$EPC_{aq} = EPC_{sed}/K_d$$

The bioconcentration factor is another measure of fate, as it reflects the chemical-specific degree of uptake by organisms from water, osmotic exchange, and equilibrium. Bioconcentration factor values and K_d values for inorganic COPC were obtained from the literature (EPA, 1986 and Baes et al., 1984).

4.2 Ecological Receptors

Evaluating risks posed by COPC to each and every species or population present is not feasible, and thus an assessment must focus on a limited number of receptors. This subset of potential ecological receptors may include organisms that are (1) chronically exposed to site-related COPC, (2) endangered, threatened, special concern, or protected species, (3) of relevance to assessment endpoints, or (4) chronically exposed via a pathway that is different from previously considered organisms.

Following is a discussion of selected receptors and the feeding habits and characteristics used to place them in the site-specific food web (Figure 4-1). These organisms were chosen based on their presence or suspected presence at the site, their trophic level, and their occupation of

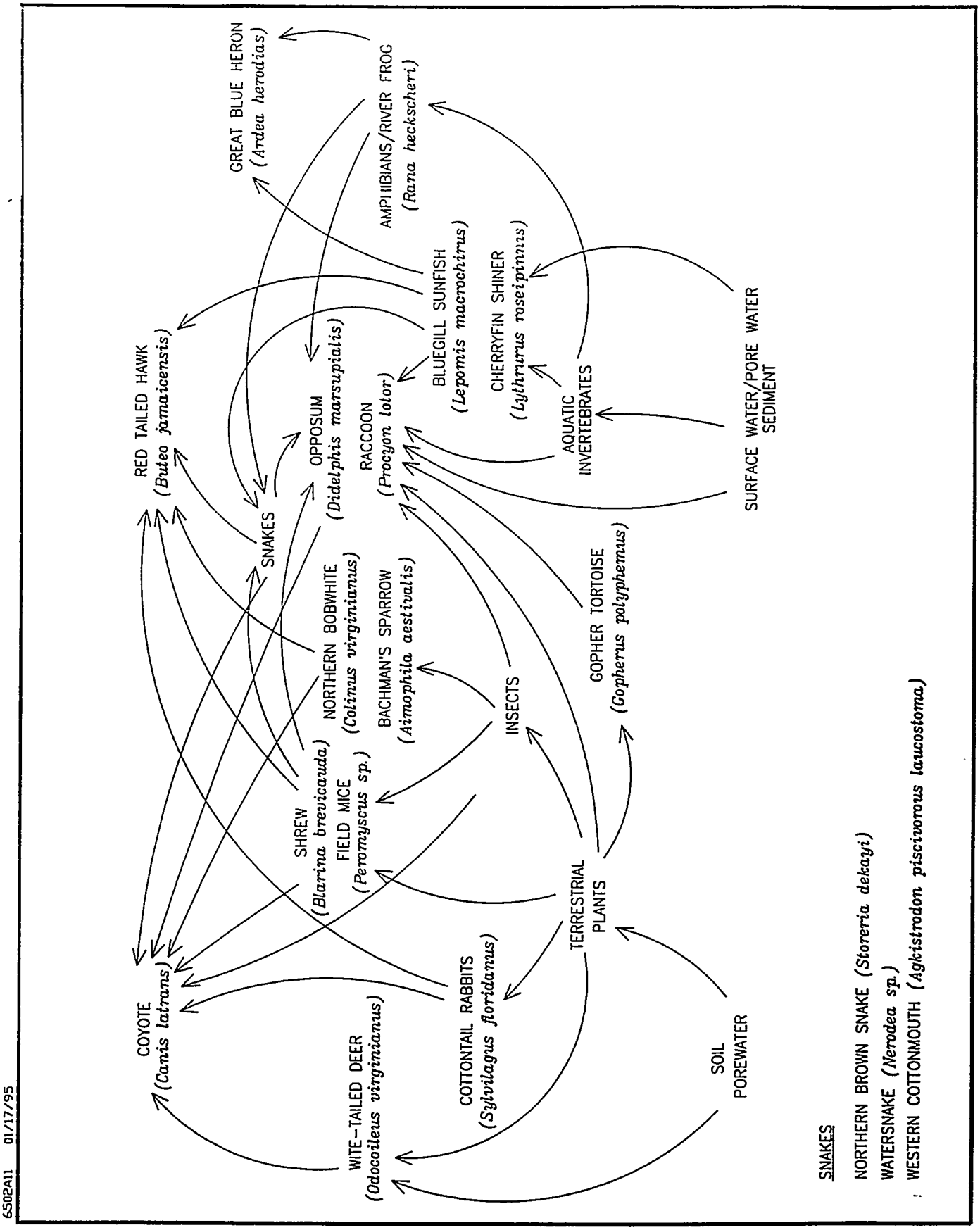

Figure 4-1
Food Web Model - Salmon Site

a specific niche not filled by another species. The chosen organisms included those sampled at the SS site.

Raccoon (Procyon lotor)

The raccoon is the most abundant and widespread medium-size omnivore in North America, except at the higher elevations of the Rocky Mountains and into southern Canada. Racoons are found near practically every aquatic habitat, particularly swamps, mangroves, floodplain forests, and freshwater and saltwater marshes. They are also common in suburban residential areas and abandoned farmlands. The raccoon is an omnivorous, primarily nocturnal feeder, eating fleshy fruits, nuts, acorns and corn, but also grains, insects, frogs, crayfish, eggs, and virtually any animal and vegetable matter (EPA, 1993). In residential areas, raccoons tip over or climb into garbage cans (Whitaker, 1993). The size of a raccoon's home range depends on its sex and age, food sources, and the season. Home ranges of a few hundred hectares appear to be most common, seasonally ranging from 5 to 5,000 hectares (2 to 2,024 acres) (EPA, 1993).

Gopher Tortoise (Gopherus polyphemus)

The gopher tortoise is a sensitive species and is the "gopher" of the Deep South. It is an accomplished burrower, creating tunnels as long as 10 m (33 ft), which slope downward from the surface, leveling off underground. The tortoise emerges daily in warm weather, usually in the cooler morning hours to forage on grass, leaves, wild fruits, and berries (Conant and Collins, 1991). No information was available on body weight, but a similar species, the Eastern Box Turtle, has a mean body weight of approximately 370 to 400 grams (EPA, 1993).

White-Tailed Deer (Odocoileus virginianus)

Once nearly exterminated in much of the Northeast and Midwest, white-tail deer are now the most plentiful game animal in eastern North America. They are found in farmland, brushy areas, and woods in most of the United States, except for a few western locations. White-tails graze on green plants and aquatic plants in the summer, and acorns, beechnuts and other nuts and corn in the fall, and feed on woody vegetation, including the twigs and buds of viburnum, birch, maple, and many conifers in the winter. Bucks and doe herd separately most of the year, but in winter gather together or "herd up" (Whitaker, 1993).

Red-Tailed Hawk (Buteo jamaicensis)

This is one of the most common and widespread members of the genus *Buteo* in the continental United States and Canada (Brown and Amadon, 1968). Red-tailed hawks live in a variety of habitats, such as farmland, woodlands, mountains, and deserts, as long as there is open country

interspersed with woods, bluffs, or streamside trees. Small mammals, including mice, shrews, voles, rabbits, and squirrels are important prey, particularly during winter. Red-tails also eat birds, lizards, snakes, and large insects, depending on availability (EPA, 1993). Home range has been reported as approximately 165 hectares (66.8 acres) (Janes. 1984). Body weights average 1,126 grams (Dunning, 1993).

Bachman's Sparrow (Aimophila aestivalis)

Bachman's sparrow is considered a sensitive species in the United States. It dwells in open stretches of pines with grass and scattered shrubs for ground cover. Like many other sparrows, it feeds on insects, such as crickets and beetles, and on seeds of grasses and sedges. It spends most of its time feeding on the ground (Whitaker, 1993). Body weights average 19.7 grams (Dunning, 1993).

Mice (Peromyscus sp.)

Mice of the genus *Peromyscus* (deer [*P. maniculatus]* or white-footed mice [*P. leucopus]*) have adapted to every possible North American habitat where they are often the most abundant mammals present (Whitaker, 1993). The habitat for mice includes all types of dry-land habitats within their range, including coniferous and deciduous forest and grasslands, as well as deserts and alpine tundra. Deer mice are omnivorous and highly opportunistic, which leads to substantial regional and seasonal variation in their diet. They eat principally seeds, arthropods, some green vegetation, roots, fruits, and fungi, as available. The nonseed plant materials provide a significant portion of the mouse's daily water requirements. Body weights of deer mice range from 15 to 35 grams, averaging 21 grams. Home range sizes vary seasonally and regionally from 0.014 to 0.128 hectares (0.006 to 0.05 acres) (EPA, 1993).

Great Blue Heron (Ardea herodias)

Although not observed at the site at the time of the species survey, this bird is likely to use the aquatic habitat at the SS and would represent a top consumer in the aquatic food chain. The heron feeds primarily on aquatic animals, preferring fish, but will also consume amphibians, reptiles, and crustaceans. Its feeding territory changes seasonally from 0.6 to 8.4 hectares (0.24 to 20.7 acres). Body weight averages 2,229 grams (EPA, 1993).

4.3 Estimated Receptor Exposures

4.3.1 Aquatic Biota Tissue Analysis

From February 26 to 28, 1993, fish tissues were collected from surface water bodies on the SS and the Reference Pond site. Sample collection was attempted at all 15 surface water stations. Although several attempts were made, no fish were obtained from Hickory Hollow Creek, Gator Pond, or Half Moon Creek Overflow Pond. Six fish samples were collected for analysis from the remaining stations. The following five species of fish were collected for whole body analysis: channel catfish (*Ictalurus punctatus*), bluegill (*Lepomis macrochirus*), starhead topminnow (*Fundulus notti*), golden shiner (*Notropis* sp.), and cherryfin shiner (*Lythrurus roseipinnis*).

Fish collection methods included dip-netting, baited hooks on trap lines, and baited killi-pots. Bait included raccoon meat, sardines, and crawfish. Due to the minimal volume of fish collected at Half Moon and Grantham Creeks, a composite sample of the species collected throughout those creeks was analyzed. An insufficient amount of fish tissue was obtained at Grantham Creek (GrC-composite), Beaver Pond-1A (BeP-1A), and Beaver Pond-2 (BeP-2) for the following analyses: alpha-total, beta-total, pesticides, and metals. However, there were sufficient fish tissues for the gamma spectrometry analysis to be performed. No analyses were carried out for cyanide or volatile or semivolatile organics.

Tissue analysis data are presented in Appendix A of this report (Tables A-2.1 through A-2.5) for inorganic, pesticide/PCBs, gross alpha, gross beta, gamma spectroscopy, and tritium analyses. All concentrations are reported on a wet weight basis.

Of the COPC for water and sediment, aluminum, barium, manganese, and zinc were found in all fish tissue samples (Reference Pond, Beaver Pond, and Half Moon Creek). Lead was found in a site sample and the reference sample at estimated concentrations below the CRDL.

No pesticides or PCBs were detected in fish samples. Gross alpha was not detected while gross beta was detected in all samples. On-site fish samples had concentrations of 0.6 and 0.9 pCi/g, and reference samples had 1.3 pCi/g. Tritium was detected in Grantham Creek, Half Moon Creek, and the Beaver Pond samples at levels ranging from 0.24 pCi/g to 0.55 pCi/g. Tritium was not detected in the Reference Pond sample. Of the other radionuclides, only potassium-40 was detected once in the Beaver Pond (7.2 pCi/g) and in the Reference Pond (2.8 pCi/g). Fish tissue detected concentrations are summarized in Table 4-1.

Table 4-1
Summary of Constituent Concentrations in Fish Tissue, Salmon Site
(Page 1 of 2)

Parameter	Reference Station	Salmon Site Stations				
		Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
METALS (mg/kg)[b] Wet weight basis						
Aluminum	13.4	2	2	3.80	2.3 BJ	5.3 B[c]J[d]
Barium	2.2 BJ	2	2	36.15	6.1 BJ	66.2
Calcium	3690.0	2	2	21,310.00	8,020.0	34,600.0
Chromium	0.35 BJ	2	2	0.32	0.25 BJ	0.38 BJ
Copper	ND[e]	2	2	2.50	1.4	3.6
Iron	41.6	2	2	18.65	13.0	24.3
Lead	0.08 BW[f]J	2	1	0.12	0.12 BJ	0.12 BJ
Magnesium	291.0	2	2	516.00	351.0	681.0
Manganese	13.4	2	2	28.65	28.2	29.1
Mercury	ND	2	1	0.43	0.43	0.43
Potassium	2910.0	2	2	2,960.00	2,790.0	3,130.0
Selenium	ND	2	1	0.34	0.345	0.345
Sodium	750.0	2	2	1,096.00	942.0	1,250.0
Thallium	ND	2	1	0.05	0.05 BWJ	0.05 BWJ
Zinc	17.6	2	2	17.35	15.1	19.6

Refer to footnotes at end of table.

Table 4-1
Summary of Constituent Concentrations in Fish Tissue, Salmon Site
(Page 2 of 2)

Parameter	Reference Station	Salmon Site Stations				
		Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
GROSS BETA (pCi/g)[g] Wet weight basis Beta	1.3	2	2	0.75	0.6	0.9
TRITIUM (pCi/g) Tritium	ND	5	3	0.43	0.24	0.55
GAMMA SPECTROSCOPY (pCi/g) Potassium-40	2.8	5	1	7.20	7.2	7.2

[a] Average concentration was calculated from detected concentrations only.
[b] Milligram per kilogram
[c] The reported value is below the Contract Required Detection Limit, but above the Instrument Detection Limit.
[d] Reported result is quantitatively estimated.
[e] Not detected
[f] Postdigestion spike for furnace atomic absorption analysis is out of control limit.
[g] Picocurie per gram

4.3.2 Terrestrial Vegetation Tissue Analysis

In conjunction with the April 26 to 30, 1993, soil sampling event, samples of terrestrial vegetation were collected for chemical and radiological analyses. The two most abundant species of pine encountered at the SS were longleaf (*Pinus palustris*) and loblolly (*Pinus taeda*), and the samples for the most part consisted of these two species. Trees sampled were on average greater than 20 years of age. The field adjacent to the catfish pond was utilized as the reference location for vegetation samples.

Vegetation samples were taken from each of the individual sites identified in Table 3-8 and Figure 3-2. Nineteen samples were taken: 18 sites plus 1 reference site. The reference sample was collected off site in the vicinity of the reference pond. Each tree sampled was cored using a ¼-in.-increment corer in order to determine the age of the specimen. This core was then used as part of the sample. The remainder of the sample mass consisted of branch segments cut into small pieces using a hand saw and pruning shears. Each of the samples was a composite formed from cores and tree branch segments taken from at least 10 trees per site. The trees sampled were no further than 60 m (200 ft) from the center of the site. Cores and clipped-branch segments were packaged in sealable plastic bags and glass jars, immediately placed on ice in coolers, and shipped to the ITAS Laboratory for analysis. Vegetation samples submitted to the laboratory were analyzed for inorganics, pesticides/PCBs, and radionuclides. These samples were not analyzed for tritium or cyanide.

The only sample that was not comprised of pine tree wood was taken from SA-6, the Helicopter Pad. No trees were within 60 m (200 ft) of the center of the sampling site; therefore, the vegetation sampled consisted of hay and a low vine that was growing in the open pasture. This sample was packaged and handled in the same fashion as the other vegetation samples.

Plant tissue analysis data are presented in Appendix A of this report. Tables A-6.1 through A-6.5 present the inorganics, pesticides and PCBs, semivolatiles, gross alpha and beta, and gamma spectroscopy data sets, respectively. The majority of inorganic constituents detected in vegetation are naturally occurring soil constituents. Many are essential nutrients. Concentrations of inorganics in vegetation on the SS were, for the most part, similar to those from the reference area. Of the COPC for soil, copper was only detected in plants from the Clean Burn Pit in SA-2 and the Helicopter Pad in SA-6 (6.70 and 6.30 mg/kg, respectively) and at the reference area (3.20 mg/kg). Lead was detected in samples from all source areas, but not in the sample from the reference area. The highest average concentration was at SA-3 (10.07 mg/kg).

Gross alpha was not detected in the reference area vegetation, but was detected in some samples from all SAs (0.26 to 3.89 pCi/g). Levels were particularly high at the helicopter landing pad in SA-6 (2.06 pCi/g) and at the South Substation in SA-4 (3.89 pCi/g). Gross beta was higher in vegetation from the reference area (1.00 pCi/g) than all other areas (0.3 to 0.9 pCi/g) except SA-6 (2.6 pCi/g). Several radionuclides were detected in the plant tissues. Cesium-137 was detected in one SA-5 sample (0.21 pCi/g) and lead-212, bismuth-214, and cerium-139 were detected only in the SA-6 sample. Tin-113 was detected only in the reference sample. Potassium-40, beryllium-7, and lead-210 were detected in samples from several locations. Table 4-2 summarizes the detected levels of constituents in vegetation on the SS and at the reference area.

4.3.3 Terrestrial Wildlife Tissue Analysis

The active collection of avian species was attempted during February 1993 using mist-netting near the junction of Main Road and Toxie Road where doves had been observed. The mist-netting was stretched across an open field in likely flight paths for 48 hours. The net was checked periodically, but the attempts proved unsuccessful. The same area was then flushed three to four times at dawn and dusk each day in hopes of being able to shoot game bird species. However, because of the time of year, game birds were scarce, and none was collected.

During the February 1993 investigation, large and small mammals were collected for metals and radionuclide tissue analysis. White-tailed deer (*Odocoileus virginianus*), cottontail rabbits (*Sylvilagus floridanus*), and marsh rabbits (*Sylvilagus palustris*) were collected with firearms under the supervision of the Mississippi Department of Game, Fish, and Wildlife. These collection activities took place on the nights of February 25 and 26, 1993. In all, seven whitetail deer and seven rabbits were collected during the two nights. Specific tissues analyzed from these specimens included fat, muscle, bone, and liver. Tissues were harvested as close to the time of death as possible, packed in appropriate containers on ice in coolers, and shipped to the ITAS Laboratory for analysis.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 1 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
METALS (mg/kg)[b] Wet weight basis	Reference	1	1	207.00	207.0	207.0
Aluminum	SA-1	4	4	107.63	61.6	159.0
	SA-2	2	2	147.00	115.0	183.0
	SA-3	4	4	100.05	87.6	115.0
	SA-4	5	5	100.12	47.6	131.0
	SA-5	2	2	201.50	159.0	244.0
	SA-6	1	1	238.00	238.0	238.0
	Sitewide	18	18	126.12	126.12	244.0
Barium	Reference	1	1	8.00	8.0 B[c]J[d]	8.0 BJ
	SA-1	4	4	11.20	2.8 BJ	16.5 BJ
	SA-2	2	2	6.75	6.4 BJ	7.1 BJ
	SA-3	4	4	7.00	6.2 BJ	7.6 BJ
	SA-4	5	5	5.64	3.4 BJ	9.0 BJ
	SA-5	2	2	12.60	8.0 BJ	17.2 BJ
	SA-6	1	1	169.00	169.0	169.0
	Sitewide	18	18	17.15	2.8 BJ	169.0
Calcium	Reference	1	1	1,650.00	1,650.0	1,650.0
	SA-1	4	4	2,002.90	1,610.0	2,430.0
	SA-2	2	2	1,591.50	923.0 BJ	2,260.0
	SA-3	4	4	1,755.00	1,260.0	2,270.0
	SA-4	5	5	2,150.00	1,470.0	3,280.0
	SA-5	2	2	2,515.00	2,330.0	2,700.0
	SA-6	1	1	4,840.00	4,840.0	4,840.0
	Sitewide	18	18	2,157.39	923.0 BJ	4,840.0

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 2 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Chromium	Reference	1	0	ND[e]	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	1	5.90	5.90	5.90
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	0	ND	ND	ND
	SA-6	1	1	0.99	0.99 BJ	0.99 BJ
	Sitewide	18	2	3.45	0.99 BJ	5.90
Copper	Reference	1	1	3.20	3.2 BJ	3.2 BJ
	SA-1	4	0	ND	ND	ND
	SA-2	2	1	6.70	6.7	6.7
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	0	ND	ND	ND
	SA-6	1	1	6.30	6.3	6.3
	Sitewide	18	2	6.50	6.3	6.7
Iron	Reference	1	1	28.90	28.9	28.9
	SA-1	4	4	25.68	16.0 BJ	37.3
	SA-2	2	2	31.30	13.9 BJ	48.7
	SA-3	4	4	21.65	16.3 BJ	25.8
	SA-4	5	5	20.24	13.2 BJ	29.8
	SA-5	2	2	24.40	20.7	28.1
	SA-6	1	1	165.00	165.0	165.0
	Sitewide	18	18	31.49	13.2 BJ	165.0

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 3 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Lead	Reference	1	0	ND	ND	ND
	SA-1	4	4	2.40	1.30	3.70 S[f]
	SA-2	2	2	2.45	1.60	3.30
	SA-3	4	4	10.07	1.60	14.30
	SA-4	5	1	4.30	4.30 S	4.30 S
	SA-5	2	1	1.90	1.90	1.90
	SA-6	1	1	2.70	2.70	2.70
	Sitewide	18	13	4.25	1.30	14.30
Magnesium	Reference	1	1	289.00	289.0 BJ	289.0 BJ
	SA-1	4	4	312.50	294.0 BJ	328.0 BJ
	SA-2	2	2	311.50	238.0 BJ	385.0 BJ
	SA-3	4	4	249.25	225.0 BJ	281.0 BJ
	SA-4	5	5	264.20	217.0 BJ	341.0 BJ
	SA-5	2	2	373.00	352.0 BJ	394.0 BJ
	SA-6	1	1	770.00	770.0 BJ	770.0 BJ
	Sitewide	18	18	317.06	217.0 BJ	770.0 BJ
Manganese	Reference	1	1	58.90	58.9	58.9
	SA-1	4	4	125.93	63.7	155.0
	SA-2	2	2	205.50	194.0	217.0
	SA-3	4	4	131.98	87.9	168.0
	SA-4	5	5	93.30	33.4	115.0
	SA-5	2	2	140.00	102.0	178.0
	SA-6	1	1	300.00	300.0	300.0
	Sitewide	18	18	138.28	33.4	300.0

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 4 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Potassium	Reference	1	1	1,010.00	1,010.0	1,010.0
	SA-1	4	3	842.67	676.0 BJ	1,090.0
	SA-2	2	2	778.50	647.0 BJ	910.0 BJ
	SA-3	4	2	680.00	606.0 BJ	754.0 BJ
	SA-4	5	5	769.40	623.0 BJ	1,000.0 BJ
	SA-5	2	2	760.50	701.0 BJ	820.0 BJ
	SA-6	1	1	623.00	623.0 BJ	623.0 BJ
	Sitewide	18	15	762.40	606.0 BJ	1,090.0 BJ
Sodium	Reference	1	1	38.20	38.2 BJ	38.2 BJ
	SA-1	4	4	50.56	42.7 BJ	65.1 BJ
	SA-2	2	2	45.00	41.6 BJ	48.4 BJ
	SA-3	4	4	52.90	37.1 BJ	63.3 BJ
	SA-4	5	5	53.60	36.1 BJ	64.2 BJ
	SA-5	2	2	50.70	48.5 BJ	52.9 BJ
	SA-6	1	1	52.50	52.5 BJ	52.5 BJ
	Sitewide	18	18	51.43	36.1 BJ	65.1 BJ
Zinc	Reference	1	1	12.20	12.2 E^{9}J	12.2 EJ
	SA-1	4	4	21.35	13.2 EJ	29.5 EJ
	SA-2	2	2	23.30	19.5 EJ	27.1 EJ
	SA-3	4	4	12.20	8.4 EJ	16.5 EJ
	SA-4	5	5	39.12	8.6 EJ	115.0 EJ
	SA-5	2	2	56.50	11.0 EJ	102.0 EJ
	SA-6	1	1	43.50	43.5 EJ	43.5 EJ
	Sitewide	18	18	29.58	8.4 EJ	115.0 EJ

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 5 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
PESTICIDES/PCBs (µg/kg)[h]	Reference	1	0	ND	ND	ND
Wet weight basis	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
Beta-BHC	SA-3	4	0	ND	ND	ND
	SA-4	5	1	27.00	27.0 J	27.0 J
	SA-5	2	0	ND	ND	ND
	SA-6	1	0	ND	ND	ND
	Sitewide	18	1	27.00	27.0 J	27.0 J
4',4-DDD	Reference	1	0	ND[i]	ND	ND
	SA-1	4	1	50.00	50.0 XHJ	50.0 XJ
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	1	88.0	88.0 XJ	88.0 XJ
	SA-6	1	0	ND	ND	ND
	Sitewide	18	2	69.00	50.0 XJ	88.0 XJ
Endrin	Reference	1	0	ND	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	2	46.00	40.0 J	52.0 J
	SA-5	2	1	42.00	42.0 J	42.0 J
	SA-6	1	0	ND	ND	ND
	Sitewide	18	3	46.00	40.0 J	52.0 J

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 6 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Endosulfan Sulfate	Reference	1	0	ND	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	1	42.00	42.0 J	42.0 J
	SA-3	4	3	38.67	35.0 J	42.0 J
	SA-4	5	3	51.67	39.0 J	71.0 J
	SA-5	2	1	49.00	49.0 J	49.0 J
	SA-6	1	0	ND	ND	ND
	Sitewide	18	8	45.25	35.0 J	71.0 J
Endrin Ketone	Reference	1	0	ND	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	4	89.50	70.0 J	130.0 J
	SA-5	2	1	87.00	87 0 J	87.0 J
	SA-6	1	0	ND	ND	ND
	Sitewide	18	5	89.00	70.0 J	130.0 J
GROSS ALPHA/BETA (pCi/g)[j]	Reference	1	0	ND	ND	ND
Wet weight basis	SA-1	4	3	0.35	0.29	0.42
	SA-2	2	2	0.33	0.26	0.40
Alpha	SA-3	4	4	0.34	0.28	0.42
	SA-4	5	3	1.55	0.39	3.89
	SA-5	2	1	0.36	0.36	0.36
	SA-6	1	1	2.06	2.06	2.06
	Sitewide	18	14	0.72	0.26	3.89

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 7 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Beta	Reference	1	1	1.00	1.0	1.0
	SA-1	4	4	0.43	0.3	0.5
	SA-2	2	2	0.40	0.3	0.5
	SA-3	4	3	0.47	0.4	0.6
	SA-4	5	4	0.63	0.4	0.9
	SA-5	2	2	0.40	0.3	0.5
	SA-6	1	1	2.60	2.6	2.6
	Sitewide	18	16	0.49	0.3	2.6
GAMMA SPECTROSCOPY (pCi/g) Wet weight basis Cesium-137	Reference	1	0	ND	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	1	0.21	0.21	0.21
	SA-6	1	0	ND	ND	ND
	Sitewide	18	1	0.21	0.21	0.21
Potassium-40	Reference	1	0	ND	ND	ND
	SA-1	4	2	6.35	6.3	6.4
	SA-2	2	1	8.60	8.6	8.6
	SA-3	4	0	ND	ND	ND
	SA-4	5	1	21.30	21.3	21.3
	SA-5	2	0	ND	ND	ND
	SA-6	1	0	ND	ND	ND
	Sitewide	18	4	10.65	6.3	21.3

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 8 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Lead-212	Reference	1	0	ND	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	0	ND	ND	ND
	SA-6	1	1	0.50	0.5	0.5
	Sitewide	18	1	0.50	0.5	0.5
Beryllium-7	Reference	1	0	ND	ND	ND
	SA-1	4	3	3.23	3.0	3.5
	SA-2	2	1	2.40	2.4	2.4
	SA-3	4	1	1.90	1.9	1.9
	SA-4	5	1	5.80	5.8	5.8
	SA-5	2	2	3.20	2.8	3.6
	SA-6	1	1	25.50	25.5	25.5
	Sitewide	18	9	5.74	1.9	25.5
Bismuth-214	Reference	1	0	ND	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	0	ND	ND	ND
	SA-6	1	1	0.68	0.68	0.68
	Sitewide	18	1	0.68	0.68	0.68

Refer to footnotes at end of table.

Table 4-2
Summary of Constituent Concentrations in Vegetation, Salmon Site
(Page 9 of 9)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Cerium-139	Reference	1	0	ND	ND	ND
	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	0	ND	ND	ND
	SA-6	1	1	0.15	0.15	0.15
	Sitewide	18	1	0.15	0.15	0.15
Tin-113	Reference	1	1	0.20	0.2	0.2
	SA-1	4	0	ND	ND	ND
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	0	ND	ND	ND
	SA-5	2	0	ND	ND	ND
	SA-6	1	0	ND	ND	ND
	Sitewide	18	1	0.20	0.2	0.2
Lead-210	Reference	1	1	2.33	2.33	2.33
	SA-1	4	2	2.91	2.29	3.52
	SA-2	2	0	ND	ND	ND
	SA-3	4	0	ND	ND	ND
	SA-4	5	2	3.18	2.21	4.15
	SA-5	2	1	2.10	2.10	2.10
	SA-6	1	1	7.61	7.61	7.61
	Sitewide	18	6	3.65	2.10	7.61

[a] Average concentration was calculated from detected concentrations only.
[b] Milligram per kilogram
[c] The reported value is below the Contract Required Detection Limit, but above Instrument Detection Limit.
[d] Reported result is quantitatively estimated.
[e] Not detected
[f] The reported value was determined by Method of Standard Addition.
[g] Inductively-coupled plasma dilution (%D) is out of control limits.
[h] Microgram per kilogram
[i] Data entered manually.
[j] Picocurie per gram

Field mice, raccoon, and opossum were collected between February 22 and 28, 1993. Traps utilized for raccoons and opossum were Conibear® Models 110 and 210. These traps were set on low tree branches or stumps and baited. Traps were set and checked three times a day. Museum Special® snap traps were used to collect field mice and other members of the rodent family. One raccoon, one Virginia opossum, one pine vole, two shorttail shrews, one Eastern wood rat, and nine field mice were collected and shipped to the laboratory for analyses. Specific tissues analyzed from the raccoon and opossum specimens included fat, muscle, bone, and liver. Large mammal tissues analyzed also included bone, liver, fat, and muscle. Whole bodies of the field mice specimens were analyzed. All small mammals collected were shipped to the ITAS Laboratory for analysis. However, of the rodent samples submitted, only one mouse sample, the woodrat, and a shrew/vole composite sample were analyzed. Trapline locations are shown in Figure 4-2.

Mammal tissue analysis data are presented in Appendix A of this report. The Appendix includes data for inorganic, pesticide/PCBs, gross alpha, gross beta, gamma spectroscopy, and tritium. Tables A-7.1 through A-7.5 present small mammal data. Tables A-8.1 through A-8.5 present large mammal data.

Many inorganic constituents were found within the mammalian tissues. Copper was detected in most tissues, but primarily in liver. Copper was only detected in bones of small mammals. Lead was detected in all tissue types in large mammals, most frequently in bone, but with the highest concentrations in fat. Similar lead results were found for small mammal tissues, with the highest concentrations in bone. The composite samples had an average concentration of 0.24 mg/kg wet weight, with a maximum of 0.36 mg/kg.

Gross alpha was detected in some large and small mammals, mostly in bone (7.29 to 7.80 pCi/g). Gross beta was detected in more samples with the highest levels in bone (21.05 to 37.43 pCi/g). Tritium was detected in some large and small mammals. Highest levels were in fat (84.59 pCi/g) and liver (6.60 pCi/g). Other radionuclides were detected sporadically in some tissues of some samples. No pesticides or PCBs were detected in mammalian tissues. The detected constituent concentrations are summarized in Table 4-3 and Table 4-4. All concentrations are presented on a wet weight basis.

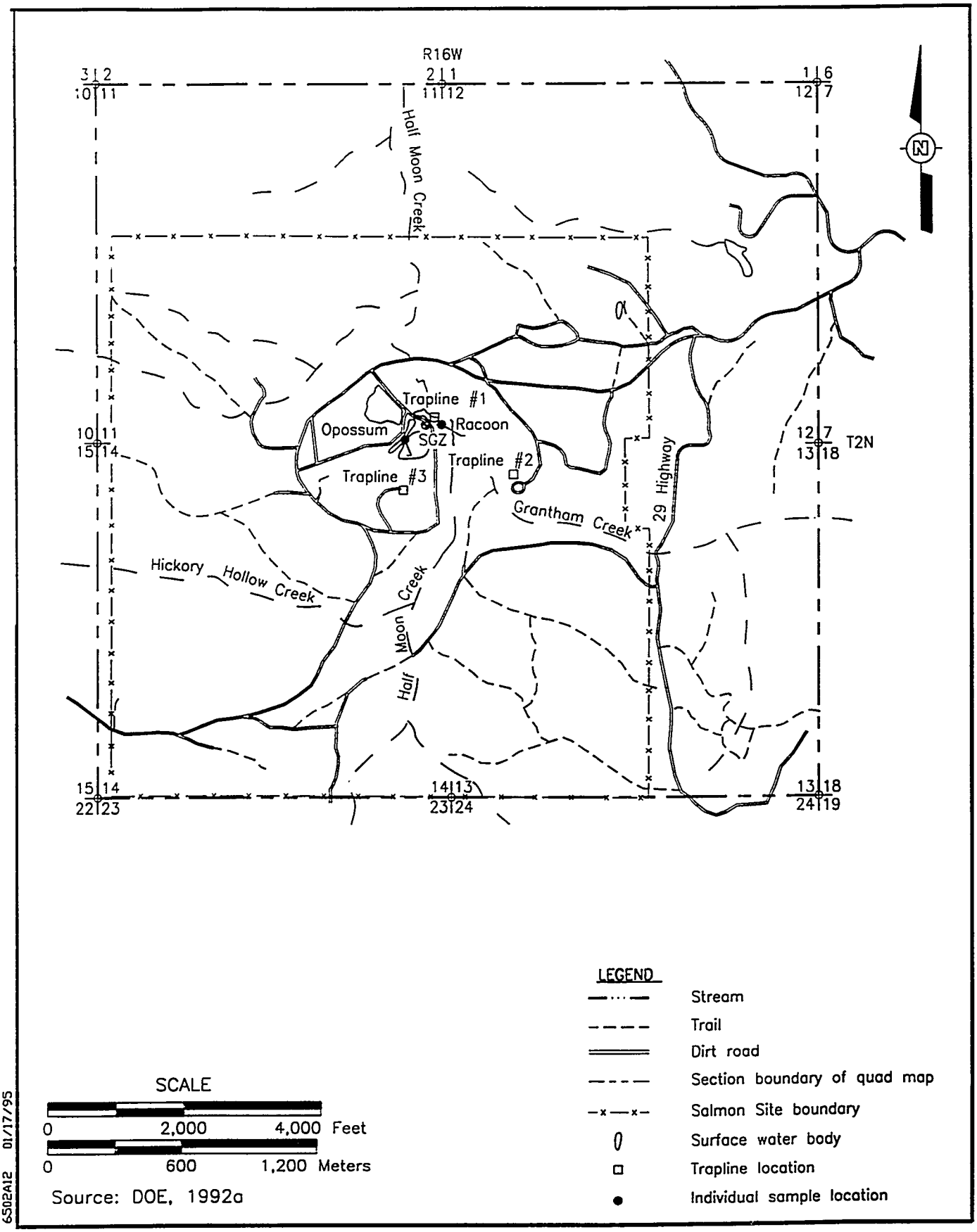

Figure 4-2
Salmon Site Location of Traplines

Table 4-3
Summary of Constituent Concentrations in Large Mammal Tissues, Salmon Site
(Page 1 of 5)

Parameter	Type	Number of Samples	Number of Detects	Average Concentration[a]	Minimum	Maximum
METALS (mg/kg)[b] Wet weight basis						
Aluminum	Bone	7	4	6.25	5.5 B[c] J[d]	7.4 BJ
	Liver	7	2	1.50	1.5 BJ	1.5 BJ
	Fat	6	2	5.55	4.3 BJ	6.8 BJ
	Muscle	7	3	1.30	1.1 BJ	1.4 BJ
Antimony	Bone	7	1	3.30	3.3	3.3
	Liver	7	0	ND[e]	ND	ND
	Fat	6	0	ND	ND	ND
	Muscle	7	2	1.25	1.1 BJ	1.4 BJ
Arsenic	Bone	7	2	1.01	0.6 2J	1.40 J
	Liver	7	0	ND	ND	ND
	Fat	6	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND
Barium	Bone	7	7	472.86	345.00	565.00
	Liver	7	7	0.46	0.12	0.94
	Fat	6	5	0.86	0.52 BJ	1.80 BJ
	Muscle	7	7	0.38	0.21 BJ	0.72 BJ
Beryllium	Bone	7	2	0.06	0.05 BJ	0.06 BJ
	Liver	7	0	ND	ND	ND
	Fat	6	1	0.22	0.22 B	0.22 B
	Muscle	7	0	ND	ND	ND
Cadmium	Bone	7	1	0.15	0.15 BJ	0.15 BJ
	Liver	7	1	0.39	0.39	0.39
	Fat	6	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND
Calcium	Bone	7	7	212,857	199,000.0	227,000
	Liver	7	7	238.79	96.5 BJ	440
	Fat	6	6	363.67	174.0 B	665
	Muscle	7	7	203.86	133.0 BJ	336

Refer to footnotes at end of table.

Table 4-3
Summary of Constituent Concentrations in Large Mammal Tissues, Salmon Site
(Page 2 of 5)

Parameter	Type	Number of Samples	Number of Detects	Average Concentration[a]	Minimum	Maximum
Chromium	Bone	7	7	3.31	2.90	3.60
	Liver	7	6	0.42	0.16 BJ	0.77 BJ
	Fat	6	5	0.47	0.26 BJ	0.71 BJ
	Muscle	7	6	0.35	0.16	0.55
Cobalt	Bone	7	1	0.29	0.29 BJ	0.29 BJ
	Liver	7	1	0.59	0.59 BJ	0.59 BJ
	Fat	6	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND
Copper	Bone	7	0	ND	ND	ND
	Liver	7	7	51.21	33.50	64.80
	Fat	6	5	2.88	0.28	5.10
	Muscle	7	7	1.79	1.50	2.20
Iron	Bone	7	6	4.87	2.8 BJ	10.4
	Liver	7	7	195.71	137	258.0
	Fat	6	4	18.33	4.0 BJ	35.8
	Muscle	7	5	29.82	24.0	35.9
Lead	Bone	7	6	2.98	0.32	13.90[f]
	Liver	7	2	3.38	0.05 BJ	6.70 S[f]
	Fat	6	3	29.62	0.27	88.20[g]
	Muscle	7	2	0.095	0.09 BJ	0.10 BW[g]J
Magnesium	Bone	7	7	3,191.43	3,050.00	3,350.00
	Liver	7	7	182.71	162.0 BJ	204.0 BJ
	Fat	6	6	63.27	25.4 BJ	91.9 BJ
	Muscle	7	7	263.14	237.0	283.0
Manganese	Bone	7	7	1.37	1.10	1.60
	Liver	7	7	5.19	4.00	6.30
	Fat	6	4	1.66	0.09 BJ	3.20
	Muscle	7	7	1.29	0.17 BJ	4.20

Refer to footnotes at end of table.

Table 4-3
Summary of Constituent Concentrations in Large Mammal Tissues, Salmon Site
(Page 3 of 5)

Parameter	Type	Number of Samples	Number of Detects	Average Concentration[a]	Minimum	Maximum
Mercury	Bone	7	0	ND	ND	ND
	Liver	7	5	0.08	0.04 BJ	0.19
	Fat	6	0	ND	ND	ND
	Muscle	7	2	0.13	0.07	0.19
Potassium	Bone	7	2	350.50	316.0	385.0
	Liver	7	7	2,672.86	2,380.0	3,200.0
	Fat	6	3	757.33	581.0 BJ	870.0 BJ
	Muscle	7	7	3,342.86	2,880.0	3,710.0
Selenium	Bone	7	2	0.62	0.57 BWJ	0.66 BE[h]J
	Liver	7	3	0.56	0.42 BWJ	0.77 BJ
	Fat	6	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND
Sodium	Bone	7	7	5,344.29	5,190.0	5,720.0
	Liver	7	7	970.29	807.0	1,260.0
	Fat	6	6	510.17	161.0 BJ	1,050.0
	Muscle	7	4	596.40	507.0	735.0
Zinc	Bone	7	7	68.53	50.7	99.1
	Liver	7	7	38.29	31.9	47.7
	Fat	6	6	4.92	2.4	7.6
	Muscle	7	7	35.26	30.3	43.7
GROSS ALPHA/BETA (pCi/g)[i] Wet weight basis Alpha	Bone	7	2	7.55	7.29	7.80 J
	Liver	7	6	0.32	0.22 J	0.40 J
	Fat	6	2	0.35	0.11	0.58 J
	Muscle	7	2	0.25	0.24 J	0.26 J
Beta	Bone	7	7	30.84	21.05	37.43
	Liver	7	7	2.61	2.27	3.06
	Fat	6	5	0.23	0.11	0.41
	Muscle	7	7	2.42	2.02	3.07

Refer to footnotes at end of table.

Table 4-3
Summary of Constituent Concentrations in Large Mammal Tissues, Salmon Site
(Page 4 of 5)

Parameter	Type	Number of Samples	Number of Detects	Average Concentration[a]	Minimum	Maximum
TRITIUM (pCi/g)	Bone	7	0	ND	ND	ND
Wet weight basis	Liver	7	2	3.36	0.11	6.60
Tritium	Fat	7	4	31.51	0.14	84.59
	Muscle	7	1	0.36	0.36	0.36
GAMMA SPECTROSCOPY (pCi/g)	Bone	7	0	ND	ND	ND
Wet weight basis	Liver	7	4	0.20	0.14	0.24
Cesium-137	Fat	7	0	ND	ND	ND
	Muscle	7	4	0.46	0.32	0.77
Potassium-40	Bone	7	0	ND	ND	ND
	Liver	7	2	6.7	3.2	10.2
	Fat	7	0	ND	ND	ND
	Muscle	7	4	3.6	2.7	4.1
Lead-212	Bone	7	4	0.28	0.00	1.00
	Liver	7	0	ND	ND	ND
	Fat	7	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND
Radium-226	Bone	7	1	0.3	0.3	0.3
	Liver	7	0	ND	ND	ND
	Fat	7	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND
Bismuth-214	Bone	7	1	0.4	0.4	0.4
	Liver	7	0	ND	ND	ND
	Fat	7	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND
Radium-228	Bone	7	1	1.40	1.40	1.40
	Liver	7	0	ND	ND	ND
	Fat	7	0	ND	ND	ND
	Muscle	7	0	ND	ND	ND

Refer to footnotes at end of table.

Table 4-3
Summary of Constituent Concentrations in Large Mammal Tissues, Salmon Site
(Page 5 of 5)

Parameter	Type	Number of Samples	Number of Detects	Average Concentration[a]	Minimum	Maximum
Titanium-208	Bone	7	0	NA	ND	ND
	Liver	7	0	NA	ND	ND
	Fat	7	1	0.30	0.30	0.30
	Muscle	7	0	ND	ND	ND

[a]Average concentration was calculated from detected concentrations only.
[b]Milligram per kilogram
[c]The reported value is below the Contract Required Detection Limit, but above Instrument Detection Limit.
[d]Reported value is quantitatively estimated.
[e]Not detected
[f]The reported value was determined by Method of Standard Addition.
[g]Postdigestion spike for furnace atomic absorption analysis is out of control limit.
[h]Inductively-coupled plasma dilution (%D) is out of control limits.
[i]Picocurie per gram

Table 4-4
Summary of Constituent Concentrations in Small Mammal Tissues, Salmon Site
(Page 1 of 5)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
METALS (mg/kg)[b] Wet weight basis	Bone	9	9	4.54	2.80 B[c] J[d]	8.00 BJ
Aluminum	Liver	1	1	1.00	1.00 BJ	1.00 BJ
	Fat	1	1	5.10	5.10 BJ	5.10 BJ
	Muscle	9	8	1.91	1.30 BJ	2.80 BJ
	Composite	3	3	10.87	4.00 BJ	17.90
Arsenic	Bone	9	1	0.75	0.75 BW[e] J	0.75 BWJ
	Liver	1	0	ND[f]	ND	ND
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	3	0	ND	ND	ND
Barium	Bone	9	9	116.39	32.10	262.00
	Liver	1	0	ND	ND	ND
	Fat	1	0	ND	ND	ND
	Muscle	9	5	0.32	0.16 BJ	0.47 B
	Composite	3	2	5.19	0.97 BJ	9.40 BJ
Calcium	Bone	9	9	57,044.44	22,500.0	98,500.0
	Liver	1	1	139.00	139.00 BJ	139.00 BJ
	Fat	1	1	251.00	251.0	251.0
	Muscle	9	9	198.00	158.0 BJ	291.0
	Composite	3	3	6,730.00	2,860.0	8,290.0
Chromium	Bone	9	9	1.63	0.33 BJ	5.10
	Liver	1	0	ND	ND	ND
	Fat	1	0	ND	ND	ND
	Muscle	9	9	0.25	0.18 BJ	0.36 BJ
	Composite	3	3	ND	ND	ND
Copper	Bone	9	7	4.69	1.20 BJ	11.10
	Liver	1	1	6.80	6.80	6.80
	Fat	1	1	5.30	5.30	5.30
	Muscle	9	9	1.61	1.10 BJ	2.20
	Composite	3	3	3.40	1.50	5.20

Refer to footnotes at end of table.

Table 4-4
Summary of Constituent Concentrations in Small Mammal Tissues, Salmon Site
(Page 2 of 5)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Iron	Bone	9	9	73.27	43.1 E^{9}J	67.5 EJ
	Liver	1	1	579.00	579.0 EJ	579.0 EJ
	Fat	1	1	30.80	30.8 EJ	30.8 EJ
	Muscle	9	9	24.89	14.2	55.2
	Composite	3	3	57.47	43.1 EJ	67.5 EJ
Lead	Bone	9	9	54.67	0.26 WJ	391.00
	Liver	1	1	0.11	0.11 BJ	0.11 BJ
	Fat	1	1	0.10	0.10 BJ	0 10 BJ
	Muscle	9	8	2.65	0.08 BWJ	19.30
	Composite	3	3	0.24	0.07 BWJ	0.36
Magnesium	Bone	9	9	1,079.11	673.0	1,530.0
	Liver	1	1	162.00	162.0 BJ	162.0 BJ
	Fat	1	1	57.30	57.3 BJ	57.3 BJ
	Muscle	9	9	291.33	250.0	319.0
	Composite	3	3	316.00	218.0	375.0
Manganese	Bone	9	9	2.18	0.38 BJ	4.60
	Liver	1	1	2.70	2.70	2.70
	Fat	1	1	0.23	0.23 BJ	0.23 BJ
	Muscle	9	8	0.53	0.21 BJ	2.00
	Composite	3	3	18.40	2.50	27.30
Mercury	Bone	9	0	ND	ND	ND
	Liver	1	1	0.64	0.64	0.64
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	3	0	ND	ND	ND
Nickel	Bone	9	2	5.30	2.40	8.20
	Liver	1	0	ND	ND	ND
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	3	0	ND	ND	ND

Refer to footnotes at end of table.

Table 4-4
Summary of Constituent Concentrations in Small Mammal Tissues, Salmon Site
(Page 3 of 5)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Potassium	Bone	9	9	2,251.11	1,340.0	2,870.0
	Liver	1	1	1,780.00	1,780.0	1,780.0
	Fat	1	1	352.00	352.0	352.0
	Muscle	9	9	3,640.00	3,210.0	3,980.0
	Composite	3	3	2,596.67	2,260.0	3,060.0
Selenium	Bone	9	1	0.07	0.07 BWJ	0.07 BWJ
	Liver	1	0	ND	ND	ND
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	3	0	ND	ND	ND
Sodium	Bone	9	9	1,693.33	1,060.0	3,080.0
	Liver	1	1	1,500.00	1,500.0	1,500.0
	Fat	1	1	759.00	759.0	759.0
	Muscle	9	9	488.11	363.0	680.0
	Composite	3	3	1,096.33	869.0	1,270.0
Vanadium	Bone	9	0	ND	ND	ND
	Liver	1	1	0.28	0.28 B	0.28 B
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	3	0	ND	ND	ND
Zinc	Bone	9	9	58.10	38.9 EJ	94.0 EJ
	Liver	1	1	25.10	25.1	25.1
	Fat	1	1	5.10	5.1	5.1
	Muscle	9	8	27.16	11.1 EJ	67.0 EJ
	Composite	3	3	22.63	18.8	25.7
GROSS ALPHA/BETA (pCi/g)[h]	Bone	9	1	12.55	12.55 J	12.55 J
Wet weight basis	Liver	1	0	ND	ND	ND
Alpha	Fat	1	0	ND	ND	ND
	Muscle	9	3	2.86	0.19	6.44
	Composite	3	0	ND	ND	ND

Refer to footnotes at end of table.

Table 4-4
Summary of Constituent Concentrations in Small Mammal Tissues, Salmon Site
(Page 4 of 5)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Beta	Bone	9	8	59.04	18.90	107.27
	Liver	1	1	1.57	1.57	1.57
	Fat	1	0	ND	ND	ND
	Muscle	9	3	2.69	1.80	3.14
	Composite	3	3	0.77	0.61	0.89
TRITIUM (pCi/g)	Bone	9	4	3.42	0.25	12.70
Wet weight basis	Liver	9	2	10.66	1.14	20.17
Tritium	Fat	1	1	0.10	0.10	0.10
	Muscle	9	2	8.88	6.37	11.38
	Composite	13	12	4.05	0.08	39.77
GAMMA SPECTROSCOPY (pCi/g)	Bone	9	2	0.21	0.14	0.28
Wet weight basis	Liver	9	3	0.50	0.31	0.60
Cesium-137	Fat	1	0	ND	ND	ND
	Muscle	9	1	0.63	0.63	0.63
	Composite	13	6	0.85	0.18	3.50
Potassium-40	Bone	9	1	2.70	2.70	2.70
	Liver	9	4	6.60	1.90	9.80
	Fat	1	0	ND	ND	ND
	Muscle	9	5	3.34	2.50	4.20
	Composite	13	5	4.26	2.80	5.60
Lead-212	Bone	9	1	0.54	0.54	0.54
	Liver	9	1	0.80	0.80	0.80
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	13	1	0.40	0.40	0.40
Uranium-238	Bone	9	0	ND	ND	ND
	Liver	9	0	ND	ND	ND
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	13	1	1.90	1.90	1.90

Refer to footnotes at end of table.

Table 4-4
Summary of Constituent Concentrations in Small Mammal Tissues, Salmon Site
(Page 5 of 5)

Parameter	Type	Number of Samples	Number of Detections	Average Concentration[a]	Minimum	Maximum
Thallium-208	Bone	9	0	ND	ND	ND
	Liver	9	2	0.14	0.13	0.14
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	13	0	ND	ND	ND
Lead-210	Bone	9	0	ND	ND	ND
	Liver	9	0	ND	ND	ND
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	13	1	1.70	1.70	1.70
Thorium-230	Bone	9	0	ND	ND	ND
	Liver	9	1	22.80	22.80	22.80
	Fat	1	0	ND	ND	ND
	Muscle	9	0	ND	ND	ND
	Composite	13	0	ND	ND	ND

[a] Average concentration was calculated from detected concentrations only.
[b] Milligram per kilogram
[c] The reported value is below the Contract Required Detection Limit, but above the Instrument Detection Limit.
[d] Reported value is quantitatively estimated.
[e] Postdigestion spike for furnace atomic absorption analysis is out of control limit.
[f] Not detected
[g] ICP dilution (%D) is out of control limits.
[h] Picocurie per gram

5.0 Ecological Effects Characterization

5.1 Potential Ecological Effects

5.1.1 Aquatic Ecosystems

5.1.1.1 Surface Waters

In order to characterize surface water COPC impacts to aquatic life, comparisons were made to the ambient water quality criteria (AWQC) for the protection of aquatic life or, when these were unavailable, to threshold values as determined from the scientific literature. The federal ambient water criteria and standards that have been established to protect the aquatic environment are the primary benchmark for comparisons. The potential for impacts to viability or propagation of aquatic life in the surface water bodies on the SS was investigated through the use of surface water concentrations for: aluminum, arsenic, and tritium (Table 5-1).

Table 5-1
Final Selection of Constituents of Concern in Surface Water, Salmon Site

Constituent of Potential Concern	BCF[a]	Surface Water Concentration (95% Upper Confidence Limit of the Mean [µg/ℓ])[b]	Chronic AWQC[c] (µg/ℓ)	Detected in Fish Tissue?	Constituent of Concern?
Aluminum	ND[d]	269.40	87	Yes[e]	Yes
Arsenic	44	0.95	48 (pentavalent) 190 (trivalent)	No	No
Iron	ND	1,813	1,000	Yes[e]	Yes
Radionuclides Exceeding Reference Surface Water Concentrations					
Tritium	NA[f]	0.34 pCi/ℓ[g]	NA	Yes	Yes[h]

[a] Bioconcentration factor, EPA, 1986
[b] Microgram per liter
[c] Ambient Water Quality Criteria
[d] No data available
[e] Detected at concentrations below the detection limit and lower than the concentrations in the reference sample. Iron was detected at concentrations above the detection limit, but lower than that of the reference sample.
[f] Not applicable
[g] Picocurie per liter
[h] All radiation sources are considered together as a constituent of concern in terms of total radiation dose.

Aluminum

Studies conducted regarding aluminum suggest that, when the pH is between 6.5 and 9.0, freshwater aquatic organisms should not be adversely affected by chronic exposure to less than 87 $\mu g/\ell$ of aluminum or by acute exposure to less than 750 $\mu g/\ell$ (EPA, 1988). Acute criteria were not exceeded in any surface water body investigated throughout the SS. However, exposure to constituents should be considered a continuous and, therefore, chronic exposure. All measured aluminum concentrations, including those from the reference pond, exceeded the chronic criteria. Therefore, it should be noted that aluminum in the surface waters at the SS are probably indicative of naturally occurring minerals in the soils and headwaters of these water bodies. Levels exceeding those from the reference pond were found at the Beaver Pond and Half Moon Creek Overflow Pond and at the furthest downstream section of Half Moon Creek. However, although aluminum was detected in fish tissue, it was at concentrations below the detection limit and lower than the concentration in the reference sample. Although unlikely, the potential for chronic toxicity to sensitive organisms exists, and aluminum is, therefore, considered a COC for freshwater systems at the SS.

Arsenic

The EPC for arsenic in surface water was 0.95 $\mu g/\ell$, which is well below the chronic AWQC of either pentavalent (48 $\mu g/\ell$) or trivalent (190 $\mu g/\ell$) arsenic. No arsenic was detected in fish tissue. Therefore, arsenic was eliminated from further considerations as a COC.

Iron

Although the 95-percent UCL of the mean for site surface water samples exceeded the reference range and the AWQC, the AWQC value was exceeded only at the Beaver Pond and Half Moon Creek overflow pond locations. However, fish-tissue samples from the Beaver Pond and Half Moon Creek had lower iron concentrations than fish from the off-site reference pond.

Radiation/Radionuclides

Tritium was the only radioactive constituent that exceeded reference concentrations in surface water. Both the Beaver Pond and Half Moon Creek Overflow Pond had tritium concentrations above the range detected at reference stations.

No apparent bioaccumulation or bioconcentration effect for tritium has been found in either aquatic or terrestrial food chains. Dilution is in fact more likely (NCRP, 1979). A major concern in tritium toxicity is undesirable genetic or cellular effects resulting from its incorporation into DNA (NCRP, 1979). However, data on levels at which these effects occur

due to tritium exposure in aquatic organisms were not available. Tritium was detected in three of five fish tissue samples, but not in the reference sample. Tritium is taken as part of the radiation dose received by aquatic organisms.

Gross beta radiation was detected in the fish tissue, but at a higher concentration in the reference than in the site samples. Alpha radiation was not detected in fish tissue. The only radionuclide detected in fish tissue was naturally-occurring potassium-40. General radiation effects are discussed in Section 5.1.2.1.

5.1.1.2 Sediment

The potential for biological effects to benthic organisms was based on COPC sediment concentrations. The assessment of effects through sediment contamination was accomplished through comparison of sediment concentrations with benchmarks.

The toxicity and kinetics of inorganics in sediment are not well understood. For this reason, risk to benthic organisms from toxic metals in sediments was characterized by comparing the measured sediment concentrations to sediment threshold values as determined from the literature. Long and Morgan (1990) have compiled data from existing studies that link concentrations of contaminants with predicted or observed biological effects. Concentrations were then selected that represent the lower tenth percentile of the screened data. This concentration is the ER-L and is used as the benchmark value in the stress characterization for metals in sediments.

COPC identified in sediment were arsenic, barium, cadmium, lead, manganese, zinc, and radiation sources, including total alpha and beta activity, tritium, and gamma spectroscopy-detected radionuclides (Table 5-2).

Arsenic

The sediment criterion derived for arsenic is 33 mg/kg (Long and Morgan, 1990). The EPC for sediment (95-percent UCL of the mean, 2.98 mg/kg) does not exceed this concentration, and arsenic would, therefore, not be expected to adversely affect aquatic life. Additionally, the calculated EPC_{aq} as an estimate of the bioavailable concentration in interstitial waters would not exceed the AWQC levels, and arsenic was not detected in fish tissues. Therefore, arsenic was eliminated from further consideration as a COC in sediment.

Table 5-2
Final Selection of Constituents of Concern in Sediment, Salmon Site

Constituents of Potential Concern	K_d[a]	BCF[b]	EPC_{aq}[c] (mg/ℓ)[d]	EPC_{sed}[e] (95% Upper Confidence Limit of the Mean [mg/kg])[f]	Sediment Criteria[g] (mg/kg)	Detected in Fish Tissue?	Constituent of Concern?
Arsenic	200	44	0.0149	2.98	33	No	No
Barium	60	N[h]	1.2923	77.54	ND	Yes	Yes
Cadmium	ND	ND	ND	2.98	5	No	No
Lead	900	49	0.0114	10.23	35	Yes[i]	No
Manganese	65	ND	3.2435	210.83	ND	Yes	Yes
Zinc	40	47	0.894	35.76	120	Yes[j]	No
Radionuclides Exceeding Reference Sediment Concentrations							
Gross Alpha	NA[k]	NA	NA	12.41 pCi/g[l]	NA	No	Yes[j,m]
Gross Beta	NA	NA	NA	16.03 pCi/g	NA	Yes	Yes[j,m]
Cesium-137	NA	NA	NA	0.40 pCi/g	NA	No	Yes[j,m]
Lead-210	NA	NA	NA	5.19 pCi/g	NA	No	Yes[j,m]
Radium-224	NA	NA	NA	1.66 pCi/g	NA	No	Yes[j,m]
Tritium	NA	NA	NA	0.16 pCi/g	NA	Yes	Yes[j,m]

[a] Soil/water partition coefficient from Baes et al., 1984
[b] Bioconcentration factor, EPA, 1986
[c] Aquatic Exposure Point Concentration. $EPC_{aq} = EPC_{sed}/K_d$
[d] Milligram per liter
[e] Sediment Exposure Concentration expressed as the 95-percent upper confidence limit of the mean
[f] Milligram per kilogram
[g] Long and Morgan, 1990
[h] No data available
[i] Analyte was detected below the Contract Required Detection Limit at estimated concentration.
[j] Average concentration did not exceed reference concentration.
[k] Not applicable
[l] Picocurie per gram
[m] All radiation sources are considered together as a Constituent of Concern in terms of total radiation dose.

Barium

No sediment criterion was available for barium. The EPA_{aq} was calculated to be 1.2923 milligrams per liter (mg/ℓ). Although barium ions are thought to be rapidly removed from solution by precipitation and other processes, concentrations in drinking water in some midwestern and western states have been known to range as high as 3 mg/ℓ (EPA, 1986). A restrictive criterion for barium does not appear warranted since barium is unlikely to exist in the toxic soluble form under usual fresh water conditions and soluble barium would have to exceed

50 mg/ℓ before toxicity to aquatic life would be expected (EPA, 1986). However, barium was detected at an elevated concentration (66.2 mg/kg) in one fish tissue sample and, therefore, warrants further consideration as a COC in sediment.

Cadmium

The ER-L for cadmium is estimated at 5.0 mg/kg (Long and Morgan, 1990). Cadmium was detected in two samples from the SS. A sample from Half Moon Creek contained 1.70 mg/kg of cadmium. One sample from the Beaver Pond (BeP-1) contained 9.00 mg/kg of cadmium. While BeP-1 exceeded the ER-L, it was found to be equal to the reported Effects Range-Median (Long and Morgan, 1990), and it is that concentration above which effects are expected to occur. Additionally, cadmium was not detected in fish tissue. Therefore, cadmium is not considered further as a COC for sediment.

Lead

The sediment criterion derived for lead is 35 mg/kg (Long and Morgan, 1990). The EPC for sediment does not exceed this concentration, and lead would, therefore, not be expected to adversely affect aquatic life at the SS. Lead was detected in one fish tissue sample, but at an estimated concentration below the CRDL. Lead is, therefore, eliminated from further consideration as a COC in sediment.

Manganese

Manganese is a micronutrient required by both plants and animals. Manganese is not considered to be a problem in fresh waters since tolerance values for aquatic life range from 1.5 to over 1,000 mg/ℓ (EPA, 1986). The calculated EPC_{aq} for manganese in SS sediments was 3.2435 mg/ℓ, at the lower end of this range. However, since the two fish-tissue samples from the site contained approximately twice as much manganese as the reference sample, manganese should be considered further as a COC in sediment.

Zinc

Zinc is a necessary micronutrient. The sediment criterion derived for zinc is 120 mg/kg (Long and Morgan, 1990). The EPC for zinc in SS sediment does not exceed this concentration and would, therefore, not be expected to adversely affect aquatic life. Zinc was detected in all fish samples, but the average concentration did not exceed the reference concentration. Zinc is, therefore, eliminated from further consideration as a COC in SS sediments.

Radiation/Radionuclides

The potential for effects exists in areas of elevated concentrations, such as the Beaver Pond and Gator Pond. Gross alpha activity in the Beaver Pond and Gator Pond were as high as 17.21 pCi/g and 22.48 pCi/g, respectively, whereas gross beta activity was as high as 18.6 pCi/g and 23.69 pCi/g. Cesium-137 was highest in the Beaver Pond (0.61 pCi/g) and the Gator Pond (1.09 pCi/g), but also occurred in the Half Moon Creek Overflow Pond, Half Moon Creek at HMC-4, and Hickory Hollow Creek at HHC-2 in levels above background. Lead-210 was elevated in the Beaver Pond as was radium-224, which was also detected at HMC-4. Tritium was only detected at the Beaver Pond. Total radiation dose was considered as a COC for sediment. See Section 5.1.2.1 for a discussion of general radiation effects.

5.1.1.3 Potential Aquatic Ecosystem Effects Summary

The final COC selected for risk characterization in the preceding sections for aquatic ecosystems at the SS include aluminum, barium, manganese, iron, and radiation sources.

The potential for chronic toxicity to sensitive organisms due to aluminum in surface waters was discussed above. Although aluminum was detected in all fish samples, the concentration was higher at the reference station than in the site samples, which were detected below the CRDL. It is likely that the aluminum concentrations in these waters are natural and that the organisms in these systems are tolerant of these concentrations. No data were available for toxicity comparisons to the fish tissue concentrations. No toxicity data were available either for manganese or barium, but it is apparent from water-quality criteria information that these elements would not be considered to be at toxic levels in SS aquatic environments. The water quality criterion for iron was exceeded in surface waters at the site, but only by the samples from the Beaver Pond and Half Moon Creek overflow pond. Fish tissue concentrations were higher in the reference pond samples than in the site samples. It is unlikely that toxicity to these organisms is occurring due to iron.

Due to the interconnected nature of the food web and the fact the terrestrial animals probably consume surface water on the site, exposure of terrestrial organisms to the aquatic contaminants should be considered. For aluminum, however, only one sample (a small mammal composite) had a concentration above the CRDL, considering all large and small mammal samples. Barium and manganese were detected in most animal tissue samples. Barium is relatively abundant in nature and tends to accumulate in the skeleton (Amdur et al., 1991). The highest concentrations at the SS were found in bone samples, and the only tissues with concentrations above the detection limit were liver samples (0.95 mg/kg wet weight) from large mammals, which may

range off site. Manganese, an essential micronutrient is present in all living organisms (Amdur et al., 1991). Manganese concentrates in the liver in the normal processing of this nutrient (Amdur et al., 1991) and was found in highest concentrations in the liver samples from the SS. These results are, therefore, likely to represent naturally-occurring concentrations. Iron is an essential nutrient also found in highest concentration in the liver samples from the SS. This element, an essential component of the oxygen-carrying molecule hemoglobin, also likely represents naturally-occurring concentrations.

Tritium, the heaviest and only radioactive isotope of hydrogen, has been a ubiquitous contaminant produced by the atomic energy programs (NCRP, 1979). However, natural tritium also exists due to cosmic ray interactions in the atmosphere (NCRP, 1979).

When tritium is introduced into an aquatic environment, the tissue water of all invertebrates and most vertebrate organisms present in that water body will approach equilibrium with the tritium concentration in the water after a period of several weeks. Fish (mosquito fish and top minnows) raised in tritiated water had tritium concentrations in their tissues approximately .31 to .56 percent of the water concentration. The tissue concentrations of tritium in fish that were fed brine shrimp, which were also raised in tritiated water, were .73 percent of the water concentration (NCRP, 1979). This was essentially the case at the SS, with a mean water concentration of 0.22 picocuries per milliliter (pCi/mℓ) and a mean tissue concentration of 0.43 pCi/g.

The radiological doses to aquatic receptors, the great blue heron and general fish species, were calculated according to the method of Baker and Soldat (1992) using the reported radionuclide concentrations in surface water and sediment. The results are shown in Table 5-3. The external dose to fish was greater than the internal dose by two orders of magnitude due primarily to the presence of cesium-137, cadmium-109, and radium-224. The internal dose was due to tritium. Tin-113 was excluded from the dose calculation due to its short half-life (115 days, Table 5-4). The heron's internal dose arises from the assumption of a 600-gram per day consumption rate of fish. The total dose for each of these receptors was well below the National Council on Radiation Protection criterion for protecting populations of aquatic organisms at 1 rad per day (NCRP, 1991) or International Atomic Energy Agency rate for terrestrial populations at 0.1 rad/day (IAEA, 1991).

Table 5-3
Total Dose due to Sediments, Surface Water, and Soil Pathways, Salmon Site

Receptor	Internal Dose (Rad/day)	External Dose (Rad/day)	Total Dose (Rad/day)
Mouse *(Peromyscus)*	$6x10^{-8}$	$6x10^{-4}$	$6x10^{-4}$
Heron *(Ardea herodias)*	$2x10^{-4a}$	$2x10^{-4b}$	$4x10^{-4}$
Fish	$1x10^{-7a}$	$1x10^{-5b}$	$1x10^{-5}$

[a]Dose due to tritium only. Dose associated with other radionuclides was not estimated due to the absence of data on the effective absorbed energy rate for the specific radionuclides from the reference document (Baker and Soldat, 1992).
[b]Greater than 50 percent of the external dose is due to cesium-137

Assumptions

1. Tin-113 was not considered when calculating dose due to the short half-life.

2. A biological half-life of one day was used when data were not available from DOE, 1994.

3. It was conservatively assumed that the probability of a gamma emission per disintegration was one in all cases. This could result in an overestimate of the external dose by approximately one order of magnitude.

Table 5-4
Selected Radionuclides of Importance and Their Half-Lives, Salmon Site

Radionuclide	Half-Life
Uranium-238	4,500,000,000 years
Plutonium-239	24,300 years
Carbon-14	5,800 years
Radium-226	1,620 years
Cesium-137	30 years
Strontium-90	28 years
Tritium	12.3 years
Krypton-85	10.9 years
Tin-113	115 days
Iodine-131	8 days
Xenon-133	5.2 days
Iodine-132	2.4 hours

5.1.2 Terrestrial Ecosystems

5.1.2.1 Soil

The assessment of potential effects to terrestrial biota, based on soil concentrations of COPC, is not as simply defined as that for surface water or sediment. For the COPC identified in soil, applied daily doses reported in the literature to cause effects were compared to soil concentrations. The COPC identified in soil are copper, lead, and radiation (Table 5-5).

Table 5-5
Final Selection of Constituents of Concern in Soil, Salmon Site

Constituents of Potential Concern	Soil Concentration (95% Upper Confidence Limit of the Mean [µg/kg])[a]	Detected in Vegetation?	Detected in Animal Tissues?	Constituent of Concern?
Copper	279.82	Yes	Yes	Yes
Lead	17.32	Yes	Yes	Yes
Radionuclides Exceeding Reference Soil Concentrations				
Lead-212	0.56 pCi/g[b]	Yes	Yes	Yes[c]
Radium-223	0.27 pCi/g	No	No	Yes[c]
Thallium-208	0.16 pCi/g	No	Yes	Yes[c]
Thorium-228	5.08 pCi/g	No	No	Yes[c]

[a]Microgram per kilogram
[b]Picocurie per gram
[c]All radiation sources are considered together as a constituent of concern in terms of total radiation dose.

Copper

Copper was detected in 34 of the 108 samples taken from the SS. The majority of the detected concentrations were within the reference range (ND - 10.70 mg/kg). One sample from Site E-2 and E-7 in SA-3 and two samples from the CH Fuel Storage Area in SA-4 had levels slightly higher than the reference area. However, all soil samples from the West Substation in SA-4 had the highest levels of copper with the minimum concentration of 25.23 mg/kg. The soil sample obtained from the center of the site activity was reported as 9,950 mg/kg, over 100 times greater than any other value. Evidence of localized copper contamination at this site is apparent.
The presence of soil COPC in vegetation tissue at concentrations above background is evidence of bioavailability and possible uptake. Copper was detected in one vegetation sample from SA-2

and one from SA-6. Copper was not detected in the vegetation samples related to the West Substation where the soil concentrations were highest.

Copper toxicity to plants may occur at very high concentrations, but it is more likely under field conditions like those at the SS that plants would become tolerant to elevated copper concentrations in soil. Copper tolerance in plants has been widely documented (Shaw, 1989).

Copper was detected in small and large mammal tissue, with highest concentrations in the liver samples. Since no reference mammal tissues were collected in conjunction with SS samples, the SS results were compared with naturally occurring reference tissue concentrations in the published literature (Table 5-6). Copper concentrations in small mammal composites were slightly above the average control area concentrations for shrew and vole species, but within the range for deer mice from one study. With only one raccoon sample and little published data, comparisons should be made with caution. The raccoon- liver copper concentration was slightly elevated compared with other omnivorous and herbivorous species. No data were available for large mammals. It is likely that these levels, being within or near reference levels, are below levels that would be expected to have an adverse effect. The tissue concentrations of copper in Bachman's sparrows and game birds are likely to be similar to that of the small mammal composite because of similar diets and possibly lower since the sparrow would have less contact with soils. The burrowing, herbivorous gopher tortoise may also have similar copper concentrations in its tissues. Copper is likely to be less toxic to the tortoise due to its slower metabolic rate.

Copper is an essential element required for enzyme processes and hemoglobin synthesis. Little information is available on chronic effects of elevated copper in biota, but may result in hemolytic anemia in humans. To determine whether bioaccumulation to higher trophic levels was a concern, copper intake was calculated for the red-tailed hawk, assuming a diet of small mammals using the following equation:

$$(I_v = \frac{(C_v)(Q_v)(FI)}{BW})$$

Table 5-6
Comparison of Copper Concentrations in Mammal Tissues to Literature Reference Values, Salmon Site

Salmon Site Species	Tissue Type	Water Content[a]	Maximum Concentration[b] (mg/kg[c] dry weight)	Literature Test Species	Tissue Type	Location	Concentration Range (mg/kg dry weight)
Small Mammal Composite	Whole Body	68%	16.25	Short-tailed shrew (*Blarina brevicauda*)	Whole Body	Not stated (U.S.)	9-11[d]
				Meadow vole (*Microtus pennsylvanicus*)	Whole Body	Not stated (U.S.)	11.9[d]
				Deer mouse Long-tailed field mouse (*Peromyscus maniculatus*)	Whole Body Whole Body	Not stated (U.S.)	10-13.4[d] 8.8-17.2[e]
Raccoon (*Procyon lotor*)	Bone	90% ash (estimated)	12.33	Raccoon (*Procyon lotor*)	Baculum bone	Illinois	0.13-1.7[e,f]
	Liver (one sample)	75% (estimated)	27.2	Rat (*Rattus rattus*)	Liver	Vermont	10.7[f]
				Rock squirrel (*Spermophilus variegatus*)	Liver	Utah	12.1-24.1[f]
				Goat (*Capra hirsus*)	Liver	Germany	11.3[f]
				Goat kid	Liver	Germany	63.3[f]

[a] EPA, 1993
[b] Salmon Site tissue concentrations are reported on wet weight basis. Values shown here are converted to a dry weight basis for comparison to literature values.
[c] Milligram per kilogram
[d] Talmage and Walton, 1991
[e] Weight on ashed basis
[f] EPA, 1980

where

I_v = intake of chemical by predator (milligrams per kilograms per day [mg/kg/d])

C_v = concentration in diet (mg/kg)

= 97.2% x 5.2 mg/kg (small mammal composite) + 2.8% x 279.82 mg/kg (soil)

Q_v = ingestion rate (kilograms per day [kg/d])

= 0.124 kg/d wet weight (EPA, 1993)

FI = fraction ingested from the site. Since the hawk's foraging area is approximately 165 hectares (67 acres) , and the site is 595 hectares (241 acres) , the calculation was made for a hawk that forages only on the SS.

= 1

BW = Body weight of hawk, kg

= 1.126 kg average for males and females, wet weight (Dunning, 1993).

$$I_v = \frac{12.89 \text{ mg/kg copper diet } (0.124 \text{ kg/d})(1)}{1.126 \text{ kg}} = 1.42 \text{ mg/kg/d}$$

It was also assumed that soil consumption would be approximately 2.8 percent using figures for the red fox (EPA, 1993), which has similar feeding habits and probably more contact with soils. An extrapolated No Observed Adverse Effect Level for copper carbonate has been calculated at 27.89 mg/kg/d and for copper oxide at 22.49 mg/kg/d for the red-tailed hawk (Opresko et al., 1993). Therefore, it is not expected that adverse effects will occur at tertiary consumer levels at SS due to copper.

Lead

Lead was detected in 105 of the 108 SS samples, ranging from 1.4 to 120 mg/kg with the 95-percent UCL of the mean being 17.32 mg/kg.

Lead was not detected in the reference area plant samples, but was detected in plants from all source areas. The highest concentration was found in a sample from SA-3 at 14.3 mg/kg. The average concentration of the detected samples sitewide was 4.25 mg/kg. Rolfe and Bazzaz (1975) saw significant uptake of lead when applying soluble lead to soil as lead chloride. Loblolly-pine and autumn-olive photosynthesis was inhibited by 11 and 17 percent, respectively, at foliar lead concentrations of 60 and 72 mg/kg dry weight, respectively. Below these concentrations, no effect was evident. Seiler and Paganelli (1987) found that red-spruce net photosynthesis was significantly lower than the control at a foliar lead concentration of only 3.2 mg/kg dry weight, but no effect was seen in loblolly pine at 2.2 mg/kg.

It is likely that foliar concentrations would be lower than or equal to the woody tissue concentrations measured at the SS. Lead tolerance in plants has also been documented and may be a factor in assessing effects to these plants. It is unlikely that lead levels seen in these plants would result in adverse effects.

Lead was detected in small mammal tissues above the CRDL in bone, muscle, and composite samples and in large mammals in bone and fat. The highest concentration was observed in a small mammal bone sample at 391 mg/kg. The SS results were compared with naturally occurring reference tissue concentrations in published literature (Table 5-7). These comparisons showed that lead concentrations in small mammals, raccoon, and deer tissues were within normal ranges and would not be expected to cause adverse effects. Again, extrapolating from these results, lead levels in the Bachman's sparrow and gopher tortoise, as well as game birds, would be expected to be in normal ranges. A calculation was made for the dose to the red-tailed hawk, using the same equation as for copper:

$$I_v = \frac{0.8349 \text{ mg copper/kg diet } (0.124 \text{ kg/d})(1)}{1.126 \text{ kg}} = 0.092 \text{ mg/kg/d}$$

C_v = Soil lead EPC of 17.32 mg/kg x 2.8% of diet + 0.36 mg/kg wet weight small mammal x 97.2%

Depression of the enzyme, aminolevulinic acid dehydratase, which is involved in hemoglobin production (and chlorophyll production in plants), occurred in all tissues of American Kestrels (*Falco sparverius*) at 25 mg/kg/d. No measurable effects were seen at 10 mg/kg/d (Eisler, 1988). The dose received here by the hawk is well below these levels, and, therefore, lead would not pose a threat of adverse effects to top consumers at the SS.

Radiation/Radionuclides

Although the 95-percent UCL of the media did not exceed the reference range, both alpha and beta activity were found above reference area levels at various sites throughout the SS. Gross alpha radiation was found at elevated levels particularly at the Clean Burn Pit and Gas Pump in SA-2, Sites E-2 and E-7 and the Government Storage Area 2 in SA-3, the Cable Storage Area and West Substation in SA-4, and the HT-2 Well Area in SA-5. Gross beta activity was found at elevated levels at many of these same sites: the Clean Burn Pit in SA-2; the CH Fuel Storage Area, Cable Storage Area, and West Substation in SA-4; and the HT-2 and HT-2M Well Areas in SA-5. Biota at these sites may be potentially affected by low levels of radioactivity, and thus radiation should be considered as a COC for soil. All radionuclides detected in site soils are

Table 5-7
Comparison of Lead Concentrations in Mammal Tissues to Literature Reference Values, Salmon Site

Salmon Site Species	Tissue Type	Water Content	Maximum Concentration[a] (mg/kg dry weight)[b]	Literature Species	Tissue Type	Location	Concentration Range (mg/kg dry weight)
Small Mammal Composite	Whole body	68%[c]	1.125	Field mouse (*Apodemus sylvaticus*)	Whole body	Not stated	1[d]
				Short-tailed shrew (*Blarina brevicauda*)	Whole body	Not stated	6[d]
				White- footed mouse (*Peramyscus leucopus*)	Whole body	Not stated	7[d]
				Deer mouse (*Peramyscus maniculatus*)	Whole body	Not stated	2.2[d]
Raccoon (*Procyon lotor*)	Liver (one sample)	75% (estimated)	0.44[e]	Raccoon (*Procyon lotor*)	Liver, kidney	Not stated	> 140[d,f]
White-Tailed Deer (*Odocoileus virginianus*)	Bone	10% (estimated)	15.44	White-tailed Deer (*Odocoileus virginianus*)	Bone	Not stated	3 - 11[d]
	Liver	75% (estimated)	26.8[g] 0.2[h]		Liver	Not stated	< 0.4[d]

[a] Salmon Site tissue concentrations are reported on wet weight basis. Values shown here are converted to a dry weight basis for comparison to literature values.
[b] Milligram per kilogram
[c] EPA, 1993
[d] Eisler, 1988. Concentrations are means unless a range is given.
[e] Based on estimated concentration below the detection limit
[f] Lead-intoxicated animal
[g] Only detection above Contract Required Detection Limit of seven samples, arrived at by Standard Addition Method
[h] Detected below detection limit at estimated concentration

considered site related and require further consideration, as discussed below. Those for which the 95-percent UCL exceeded the reference range of concentrations included lead-210, thallium-208, and thorium-228.

Vegetation in SA-6 had the highest gross alpha and beta measurements. Vegetation from SA-6 was entirely crop plants, and the samples are not directly comparable since all other samples were composed of pine. Excluding SA-6, gross beta was higher in reference area vegetation than on-site vegetation. Gross alpha was similar on site to reference areas except for SA-4. Of the COPC identified for soils, copper, alpha and beta radiation, and the radionuclides cesium-137, potassium-40, lead-212, beryllium-7, bismuth-214, cerium-139, tin-113, and lead-210 were detected in vegetation. Most radionuclides were detected in only 1 of the 18 samples, except lead-210 (in 6 samples) and the naturally occurring beryllium-7 (in 9 samples) and potassium-40 (in 4 samples). No tritium analyses were conducted for either plants or soils. Gross alpha and beta radiation and several radionuclides were detected in small and large mammal tissues. Radionuclides detected in more than 1 sample include cesium-137, potassium-40, lead-212, and thallium-208 (detected in 2 small mammals and 1 large mammal). A general discussion of radionuclide effects is as follows.

Radioactive atoms, which represent unstable forms of an element, undergo spontaneous nuclear transformations. Charged particles and electromagnetic waves are given off in the form of radiation. Excess energy released in this way is termed ionizing radiation. Such transformations are referred to as radioactive decay. A fundamental and unique characteristic of each radionuclide is its radioactive half-life. The half-life of a radionuclide is the time required for the number of atoms present to decrease (decay) by one-half. Half-lives range from fractions of a second to millions of years. Selected radionuclides of potential importance for the SS and their half-lives are listed in Table 5-4.

The absorption of the energy of ionizing radiation in cells involves ionization of atoms and the production of ions within the cells. Although the exact mechanism of action of ionizing radiation is not known, radiation injury is considered to be related in some way to the transfer of energy into a cell and the production of ions within the cell. Ionizing radiation can cause deleterious effects on biological tissues only when the energy released during radioactive decay is absorbed by the tissue. The dose delivered to tissues from radiations external to the body occurs only while the radiation field is present. However, the dose delivered to body tissues due to radiations from systemically incorporated radionuclides may continue long after intake of the nuclide has ceased.

Alpha particles originate in the nuclei of radioactive heavy elements. Because of their double positive charge, these particles have great ionizing power, but their large size results in very little penetrating power. Alpha particles typically travel only about 4 cm (1.6 in.) through air and cannot penetrate a sheet of paper or a single layer of skin. However, alpha-emitting radionuclides are considered internal radiation hazards and are dangerous if inhaled or ingested.

Beta particles are emitted from within the nuclei of atoms undergoing a radioactive decay process. Many of the common isotopes used in biological research, such as tritium, carbon-14, phosphorus-32, sulfur-35, and calcium-45, emit beta particles. The ability of beta particles to penetrate a material is dependent upon the energy of the particle. Radionuclides that emit beta particles are considered both external and internal radiation hazards. The probability and severity of risk from beta emitters is related to the kinetic energy of the beta particle, the physical half-life of the radionuclide, and the biological half-life of the radioisotope. Beta particles from tritium do not have enough energy to penetrate a single layer of dead skin cells (Shapiro, 1981).

Gamma rays originate from unstable atomic nuclei releasing energy to gain stability. They have definite energies, characteristic of the nuclide from which they are emitted. Gamma rays ionize materials largely indirectly through a variety of mechanisms that involve ejection of high-speed electrons from the atoms by which they are absorbed. Because of their high penetration ability, gamma rays are of most concern as external hazards.

The sources for exposure of plants and animals to ionizing radiation can be broken down into four major groups: (1) natural sources of irradiation, both external and internal; (2) medical sources and radiopharmaceuticals; (3) nuclear reactions, such as nuclear power reactors and nuclear weapons; and (4) other miscellaneous sources. Exposure to natural sources of irradiation are unavoidable for the most part, but the degree of exposure to man-made sources is subject to change, depending upon usage of such sources.

The debris from nuclear detonations predominantly emit gamma and beta radiation. Some of the more common radionuclides involved in a nuclear explosion are uranium-238, plutonium-239, carbon-14, radium-226, cesium-137, strontium-90, tritium, krypton-85, iodine-131, xenon-133, and iodine-132.

The principal exposure and primary dose from nonfallout radiation include the buildup of deposited radionuclides in soil and vegetation, transfer of radioactive materials from soil to plants, and incorporation of radioactive materials in animals, both prey and predatory species. Currently,

the projected annual average whole-body dose rate from global atmospheric weapons testing is projected to be 4 to 5 millirems per year through the year 2000 for the United States population (BEIR, 1980). No estimates are given for wildlife, but one could assume them to be equal to, or slightly higher than, those given for humans.

Although different radionuclides have different properties and energies, alpha, beta and gamma radiation are all sources of ionizing radiation and, since they occur together at the site, would all contribute to the total dose. Dosages are measured in rads or grays. One hundred rads are equal to one gray.

Radionuclide levels in tissues can change with dietary composition (e.g., different levels may occur in different types of vegetation; an animal may shift from vegetation to insects at different times of the year). Therefore, the radiation dose was calculated to the mouse as a terrestrial ecological receptor that is assumed to spend all of its time on site. The radiological dose to the mouse was calculated using radionuclide soil concentrations and a modification of the methodology outlined in the *Hanford Site Risk Assessment Methodology* (DOE, 1994). Yu et al. (1993) was also consulted.

The external dose to the mouse was about four orders of magnitude greater than the internal dose. The primary contributors to the external dose were cesium-137 and thorium-228. For the external exposure calculation, it was conservatively assumed that, for each radionuclide that is a gamma emitter, the probability of a gamma emission is one. This means that the mega-electron volts per disintegration used in the calculation was not multiplied by the probability of gamma being the mode of decay. The resulting external dose was less than 1 rad per day. Therefore, this conservatism is permissible. The total dose to the mouse was 6×10^{-4} rad per day (Table 5-3), well below the IAEA (1991) rate of 0.1 rad per day for terrestrial population protection. The Bachman's sparrow is expected to have a similar internal dose due to diet similarities between this sensitive species and the mouse. However, the external dose would probably be lower since the sparrow does not burrow and would spend less time in contact with the soil. The dose to the tortoise is expected to be as low as the mouse, or lower, since the external dose would be reduced by the shielding effect of the shell (Cosgrove, 1971).

Woodwell and Rebuck (1967) studied chronic gamma irradiation of an oak-pine forest. Pines were the most sensitive species, followed by shrubs, herbacious plants, and lichens. A reduction

in species diversity was observed at 100 rad per day (1 gray per day) (Woodwell and Rebuck, 1967). Measurable changes in diversity in a grassland plant community require a dose rate of about 400 rad/day, and a calculation using effective decay energy for Pu-239 was devised to determine the concentration required to achieve this dose rate (Whicker, 1980). It was determined that 1.5 x 10^{-2} microcuries (μCi) Pu-239/g plant tissue or 150 μCi Pu-239/g soil were required to achieve this dosage (Whicker, 1980). The dose to the SS plants is well below these levels.

5.1.2.2 Potential Terrestrial Ecosystem Effects Summary

Levels of lead and copper in plant and animal tissues at the SS were found to be generally within reference ranges found in the published literature and would, therefore, be below levels that would be expected to cause adverse effects. Calculations of the intake of these elements by the top predator observed at the site, the red-tailed hawk, also gave results for which adverse effects would not be expected. The radiation dose calculated for the mouse was well below established criteria for radiation exposure effects to terrestrial populations. No adverse effects to the plant and animal populations at the SS are expected due to copper, lead, or radiation sources.

5.1.3 Laboratory Studies

5.1.3.1 Aquatic Toxicity Tests

Toxicity tests were used to assess the potential for gross effects from contaminated media on selected test species. The toxicity tests integrate the acute effects of all constituents with the bioavailability of those constituents in the contaminated media.

Surface water samples used in toxicity testing were collected on April 25, 1993, from the fifteen stations indicated in Table 3-6 and 3-7. All samples were collected as composited, mid-depth samples from random points at each station. Samples were immediately packed on ice and shipped for testing as outlined in the EEW Standard Operating Procedures. Ceriodaphnid and fathead minnow tests screened surface waters for acute toxicity. The assessment of toxicity is made by comparing the response, either survival or germination, in the on-site media to the response in the reference area media.

The results of toxicity tests are presented in Tables A-4.10 and A-4.11 in Appendix A of this report. Survival of fathead minnows in surface waters of the SS was at least 95 percent in all site samples, references, and controls. Survival of ceriodaphnids in test waters was lower. The lowest survival was seen in the ponds and Hickory Hollow Creek. Survival in on-site ponds

ranged from 60 to 72 percent, but was only 80 percent in the reference pond. On-site Hickory Hollow Creek survival was 64 percent; upstream survival was only 60 percent. Grantham Creek survival ranged from 80 percent on site to 88 percent upstream. Half Moon Creek survival ranged from 88 to 100 percent on site as compared to 88 percent upstream. No significant differences were found between any on-site location and the applicable reference with respect to acute toxicity.

5.1.3.2 Terrestrial Toxicity Tests

Soil samples used in toxicity tests were collected as aliquots of samples collected for chemical and radionuclide analyses. These samples were collected from April 26 to 30, 1993. Samples were immediately packed on ice in a cooler and shipped for testing as outlined previously. Samples were not composited, but tested as individual replicates for each station. The lettuce seed germination test was used to screen surficial soils for acute toxicity. The assessment of toxicity is made by comparing the response, either survival or germination, in the on-site media to the response in the reference area media. Testing was conducted from May 4 to 9, 1993, in the IT-Edison Bioassay Laboratory. Average seed germination per site soil sample ranged from 64 percent to 72 percent. Reference area germination was 66 percent and laboratory control was 64 percent.

5.1.4 Field Investigations

5.1.4.1 Benthic Ecology Investigations (Rapid Bioassessment Protocol I)

During the period July 13 to 19, 1992, each of the nine creek stations underwent benthic macroinvertebrate analysis using RBP I (EPA, 1989c). The six pond stations were surveyed to support general macrobenthos quality comparisons.

Benthic macroinvertebrate samples were collected using a kick-net technique at all lotic system (running water) stations and with an Ekman grab sampler at all lentic-system stations. Data collected from the samples were relative abundance of macroinvertebrate families or orders, as well as quality measurements (e.g., dissolved oxygen and pH) and habitat characteristics (e.g., substrate, flow, and cover) for general habitat assessment.

The benthic-community habitat-assessment data for the ponds and creeks of the SS are presented in Appendix A. Table A-9.1 is the habitat assessment for RBP I and Table A-9.2 is the RBP I level macrobenthos data.

The RBP I habitat assessment found all creek habitats to be comparable to the reference station or of a higher quality, except for station HHC-2. This station was evaluated as capable of supporting a habitat similar to its reference station (station HHC-1).

The macrobenthos qualitative field data from the RBP I investigation are summarized in Appendix Table A-9.2. Ten taxa were observed at the various creek stations. Taxa identification and confirmation was performed using Peckarsky et al., 1990. A description of the ecological requirements and importance of the taxa found at the SS was presented in the *Results of a Preliminary Survey for Threatened and Endangered Species and Floodplains and Wetlands at the TDTS* (IT, 1992a). The most prevalent taxa found throughout the site were Plecoptera, Coleoptera, Anisoptera, Decapoda, and Tabanidae. Station HHC-2 was the lowest in taxa richness and was abundant in Chironomidae, which are generally a stress-tolerant species. This was probably more attributable to the station's creek topography than external sources of stress, due to the abundance of pollution intolerant Anisoptera that were also found at that station. Station HMC-3 had a large submerged root mass present that was found to support a diverse macroinvertebrate population. Grass shrimp and Gastropods were unique to that station, which is located closest in proximity to SGZ.

The SS aquatic ecosystem was evaluated using the Habitat Assessment/Impairment technique associated with the RBP (EPA, 1989c). The data for the RBP I are presented in Appendix Table A-9.1. The RBP Habitat Assessment compares the ratio of the score for the station of interest to the score of the appropriate reference station. The result of this comparison indicates whether the habitat across the site is capable of supporting an adequate aquatic community. The results for the RBP I habitat assessment found that HHC-2 had the lowest percent comparability to its reference station, yet it is still classified as supporting adequate biota. This may be attributed to the fact that Hickory Hollow Creek is an intermittent stream that was reduced to standing pools in the area of HHC-2 during the study period. However, the sitewide RBP I results, as presented in Appendix Table A-9.1, reflect habitat conditions and parameters that are predominantly in the good to excellent category for supporting a biological community.

The RBP I qualitative evaluation showed that stations HMC-2, HMC-5, and GrC-2 shared a common overall abundance and taxa richness, with three different taxa present. Station HHC-2, had two taxa that were abundant (>10 individuals) and three taxa that were rare (<3 individuals). Stations HHC-2, GRC-1, HMC-1, HCM-3, and HMC-4 shared a common overall abundance and

taxa richness, with four to seven different taxa occurring at those stations. Crayfish, lampreys, salamanders, tadpoles, Anisoptera, Coleoptera, and Plecoptera were found commonly throughout the creek areas. The presence of these taxa is normally indicative of an uncontaminated habitat.

5.1.4.2 Benthic Ecology Investigations (Rapid Bioassessment Protocol III)

The RBP III was used for a more quantitative benthic ecology survey and was conducted during the second phase of the investigation between February 22 and 28, 1993. The 15 stations sampled previously using RBP I were resampled using RBP III. A kick net was again employed for collecting both riffle/run and pool samples. In addition, a Coarse Particulate Organic Matter (CPOM) sample was collected from each station for functional feeding group analysis. Appendix Table A-9.3 is the habitat assessment for RBP III, Appendix Table A-9.4 is the RBP III level macrobenthos data, and Appendix Table A-9.5 is the CPOM data..

The RBP III habitat assessment found Half Moon Creek stations to be comparable to the reference station or of a higher quality. Both Grantham Creek and Hickory Hollow Creek were found to be capable of supporting a habitat similar to their respective reference stations (station HHC-1 and GrC-1).

The macrobenthos field results for the RBP III investigation are summarized in Appendix Table A-9.4. Sixty-seven taxa were collected in the riffle/run samples. The only taxa that were identified in the RBP III investigation that were not found in the RBP I investigation were Hirudinea (leeches), Turbellaria (flatworms), and Megaloptera (alderflies). All taxa were identified at the IT Aquatic Sciences Laboratory in Edison, New Jersey, to the lowest practical taxa. The most prevalent taxa found throughout the site were Plecoptera, Tricoptera, Ephemeroptera, and Diptera. The biological integrity of each creek station was evaluated by performing the analysis for the eight metrics that are associated with the RBP III.

The SS aquatic ecosystem was evaluated using the Habitat Assessment/Impairment technique associated with the RBP (EPA, 1989c). The data for the RBP III are presented in Appendix Table A-9.3. The RBP Habitat Assessment compares the ratio of the score for the station of interest to the score of the appropriate reference station. The result of this comparison indicates whether the habitat across the site is capable of supporting an adequate aquatic community. The results for the RBP III habitat assessment found that both HHC-2 and GrC-2 were capable of supporting a similar biological community as compared to their respective reference stations and that all HMC stations were comparable or of a higher quality than the reference station. The conditions at HHC-2 may be attributed to the intermittent nature of that creek, as discussed

previously. The conditions at GrC-2 are capable of supporting a biological community that is similar to the reference station as demonstrated by the results of the benthic macroinvertebrate investigation (as presented in Appendix Table A-9.5). The data demonstrate that GrC-2 had 20 organisms present, representing 11 different species. as compared to 4 organisms present that represented 4 different species at station GrC-1. The overall RBP III Habitat Assessment results, as presented in Appendix Table A-9.3, reflect habitat conditions and parameters that are predominantly in the good to excellent category for supporting a biological community.

The RBP III is a more rigorous bioassessment technique than the RBP I, involving systematic field collection and subsequent lab analysis in order to detect more subtle degrees of impairment. The RBP III macrobenthos data are summarized in Table A-9.4. The final analysis of the RBP III is a summation of biological scores from eight different metrics that evaluate the biological integrity at each creek station as compared to its respective reference station. Using the raw benthic data (see Appendix Table A-9.4), a numeric value is calculated for each metric. Calculated values are then compared to values derived from either an unimpaired reference site within the same region or a suitable control station on the same stream. The scores for the eight metrics are then totaled and compared to the total metric score for the reference station. The first seven metrics evaluate the riffle/run macrobenthos and metric #8 evaluates the functional feeding groups present in the CPOM sample. The percent comparison between the total scores provides a final evaluation of relative biological condition.

The results of the RBP III analysis are presented in Table 5-8. The range of biological scores for the reference stations ranged from 48 at HMC-1 to 42 at GrC-1. The lowest biological score of 14 of all the test stations occurred at station HMC-5 and the highest biological score of 42 occurred at station HMC-4. The percent comparability to the reference station indicated that HMC-2 and HMC-3 as being slightly impaired due to the community structure being less than expected.

Station HMC-5 was found to be moderately impaired as compared to the reference, due to a reduction in the Ephemeroptera, Plecoptera, Trichoptera index and the presence of tolerant types of species. The habitat assessment for the Half Moon Creek system was found to be comparable to the reference area, and good finfish diversity was found throughout the entire creek system. The grain-size distribution across the site and within the Half Moon Creek System was found to be uniform and to be approximately 90-percent sand. The grain-size data were further discussed in Section 3.2.1.2.

Table 5-8
Benthic Ecology RBP III Metric Results, Salmon Site

Station Location	Biological Score								Total
	Metric #								
	1	2	3	4	5	6	7	8	
GrC-1[a]	6	6	6	6	-[b]	6	6	6	42
GrC-2	6	6	-	-	0	6	6	6	30
HHC-1[a]	6	6	6	6	4	6	6	6	46
HHC-2	6	6	-	2	6	6	6	6	38
HMC-1[a]	6	6	6	6	6	6	6	6	48
HMC-2	6	6	-	2	6	6	4	6	36
HMC-3	6	6	6	2	4	2	4	2	32
HMC-4	6	6	6	4	6	6	6	2	42
HMC-5	2	6	-	2	0	0	2	2	14

Station Location	% Comparability								% Comparability to Reference	Biological Condition Category
	Metric #									
	1	2	3	4	5	6	7	8		
GrC-1[a]	100	100	100	-	-	100	-	100	100	Nonimpaired
GrC-2	155	90	0	-	0	133	-	171	92	Nonimpaired
HHC-1[a]	100	100	100	100	4	100	-	100	104	Nonimpaired
HHC-2	151	115	200	40	6	200	-	55	133	Nonimpaired
HMC-1[a]	100	100	100	100	6	100	-	100	106	Nonimpaired
HMC-2	130	95	0	25	6	100	-	80	78	Slightly
HMC-3	93	82	100	25	4	75	-	85	72	Slightly
HMC-4	128	114	200	70	6	150	-	20	120	Nonimpaired
HMC-5	43	101	0	25	0	25	-	22	36	Moderately

[a]Reference station
[b]Indicates that the metric value could not to be calculated

5.1.4.3 Aquatic Field Observations

The RBP habitat assessment technique is appropriate for streams and rivers only; therefore the ponds on site were not evaluated using this technique. The pond habitats had a predominantly muck-mud or organic-detritus substrate, including coarse and fine particulate organic matter. The inorganic substrate consisted predominantly of sand mixed with silt or clay. The canopy is open at the Beaver Pond and the Reference Pond (ReP-1), and a partly shaded canopy exists for the Gator Pond (GaP-1) and Half Moon Creek Overflow Pond (HOP-1). All ponds except for HOP-1 and ReP-1 had water lilies present on approximately 75 to 80 percent of their surface area. The pond habitats that were observed are typical for the site topography and drainage basin area.

In support of the aquatic field observations, a finfish survey was conducted during the threatened and endangered species survey (IT, 1992a). The finfish biosurvey results indicated a diverse and abundant population present for both the creek and pond stations at the SS. Sufficient food sources and suitable habitat exist for the species found across the site, which appears to further indicate a healthy aquatic ecosystem.

The SS pond stations were not evaluated using the RBP because this method is not applicable to nonlotic systems. The pond stations had similar numbers of taxa except for the Beaver Pond where a great number of diptera species were found. At station BeP-1, 291 of the 298 organisms that were found were diptera species. At station BeP-1a and BeP-2 approximately 50 percent of the organisms found were diptera species. The diptera or chironomidae species tend to be present in greater numbers in stressed aquatic systems, due to the chironomidae being generally a pollutant-tolerant species. Therefore, some biological alteration was found at the Beaver Pond as compared to the reference pond site.

5.1.4.4 Terrestrial Field Observations

Field observations of the SS terrestrial ecosystem occurred during July 13 to 19, 1992. The results of the vegetation, wetland/floodplain, avian, herpetological, and mammalian surveys conducted were presented in detail in the *Results of a Preliminary Survey for Threatened and Endangered Species and Floodplains/Wetlands at the SS* (IT, 1992a). The Phase II investigation, which occurred in February and April 1993, involved additional field observations made during the collection of soil, vegetation, and large and small mammal samples. The terrestrial ecosystem of the SS was found to have good diversity, as well as no observed abnormalities associated with the vegetation, soil, or biota.

6.0 Risk Characterization

6.1 Risk Estimation

For the assessment at the SS. the following assumptions were made:

- COC and their respective concentrations in on-site media were appropriate indicators of overall contamination.
- Bioassay tests performed with media collected from the site were appropriate indicators of gross biological effect.
- Field observations made on a qualitative basis were indicative of overall community and individual stress.
- The reference sites selected were similar to the study areas with regard to environmental characteristics. but were relatively unaffected by site-related activities and contamination.
- Only those constituents with site concentrations greater than reference area concentrations (for organics and inorganics) were considered further as possible COC.

Risks posed by COC concentrations within the surface water, sediment, soil, and biological tissues were assessed through comparisons with benchmarks. These benchmarks were criteria or standards gathered through review of the scientific literature. Contaminant concentrations exceeding benchmarks are indicative of potentially stressful conditions.

Toxicity test response data from the on-site samples and reference sample were compared and any on-site sample showing statistically significant adverse responses relative to the reference sample was considered to show measurable toxicity.

Comparisons between on-site community conditions and reference conditions or expected conditions were made using biological data gathered from benthic macroinvertebrate analyses or terrestrial site observations. In addition, chemical concentrations within biotic tissue could offer evidence of exposure, which may be detrimental or lead to food chain transfers.

6.2 Risk Description

Section 3.3 of this report described the exposure pathways and the factors relating to them. Five pathways were described that have the potential to expose biota to COCs: water, sediment, soil, air, and biota via food chain transfers. The following sections describe the potential risk to ecological receptors posed by each pathway, within the aquatic and terrestrial ecosystems.

6.2.1 Aquatic Ecosystem

6.2.1.1 Water Pathway

Some impairment of benthic communities according to RBP III metrics was detected at stations HMC-2, HMC-3, HMC-5, BeP-1, BeP-1a, and BeP-2. Impairment of benthic communities was also detected at stations HMC-5. BeP-1, BeP-1a, and BeP-2, as well as stations HMC-4, HOP-1, and GaP-1. This impairment could be due to elevated contaminant concentrations or to habitat factors. Impairment based on acute toxicity testing was not observed at any stations.

There is strong evidence for absence of constituent-induced risk at both GrC-2 and HHC-2 since impairment was not detected according to analytical data, bioassay tests, or field observations. The slight impairment at HMC-2 and HMC-3 is probably not due to constituents because the COC concentrations were not elevated with respect to the reference areas. The measured constituents in the Half Moon Creek Overflow Pond, the Gator Pond, and HMC-4, though elevated, may not be bioavailable to organisms because no impairment was found. HMC-5 had a high concentration of aluminum (235 $\mu g/\ell$). This exceeds the chronic criteria of 87 $\mu g/\ell$ by a factor of almost three. Although no acute toxicity was found in these surface waters, elevated aluminum may be causing chronic risk to organisms and, thus, may be the cause of the slight impairment in benthic macroinvertebrate communities there, according to RBP III metric interpretation. Risk to the fish population due to aluminum or iron would appear to be negligible because the site tissue samples had lower concentrations than the reference sample.

Tritium contributed the entire internal dose to fish, but the total dose, including that received from sediments, was still well below the protective criterion for aquatic populations. The dose to the great blue heron was also extremely low and within protective dose rates. Radiation thus poses no risk to aquatic biota through the water pathway.

6.2.1.2 Sediment Pathway

The Beaver Pond had high levels of many constituents. Although surface waters were not found to be acutely toxic, the benthic macroinvertebrate community appeared stressed relative to the reference area. The elevated levels of COC in the sediments may be causing chronic effects on aquatic biota in the Beaver Pond.

On an over-all basis, the only apparent potential for risk would be related to the presence of elevated levels of barium and manganese. However, although site fish tissues had elevated concentrations of these elements compared to the reference sample, adverse effects would not be

expected in the aquatic life on the site because neither barium nor manganese exists in high soluble (bioavailable) concentrations in fresh water. Neither chemical was considered a COPC in surface water. Because these two metals are widely found in nature and manganese is a micronutrient, it is likely that the concentrations in the fish are within normal ranges and do not represent a sitewide risk. The lack of effects data for manganese and barium in fish tissue, however, could be considered a data gap.

Radiation, as discussed above, was calculated as a total dose to fish and herons from water and sediment sources and poses no risk to aquatic biota.

6.2.2 Terrestrial Ecosystems

6.2.2.1 Soil Pathway

Comparison of SS tissue results for lead and copper to literature reference values and calculations of dose to a top predator led to the conclusion that no adverse effects would be expected for terrestrial receptors on a sitewide basis. The same is true for radiation sources, for which calculated dose to the mouse was well below the protective criterion for terrestrial populations. It is reasonable to extrapolate these results to the Bachman's sparrow, game birds, and gopher tortoise because of the similarity in their diets. The birds would have less contact with soils than the mouse, and although the gopher tortoise may have more soil contact, its shell is an effective shield against radiation (Cosgrove, 1971).

There were no observed signs of stress at any of the sites except for the physical alterations due to DOE-related activities. Also, based on laboratory toxicity tests, there was no evidence of acute toxicity to flora in the surficial soils. However, COC were present at a few sites. The Clean Burn Pit in SA-2 and the HT-2 Well Area in SA-5 had high levels of both alpha and beta activity and some radionuclides. The West Substation in SA-4 also had high measurements of gross alpha and beta activity and radionuclides and also had high levels of copper. Although the potential for risk to terrestrial receptor populations site-wide is negligible, there may be risk of chronic effects to individuals in these localized areas.

6.2.2.2 Air Pathway

There is no ecological risk expected related to the air pathway. The only release mechanism to air for lead and copper would be through adsorption to soil-particles and mobilization to air as fugitive dust. A similar mechanism is also possible for radionuclides, and minor levels of noble gases, such as krypton and argon, may have been emitted during reentry drilling operations.

However, as discussed in Section 3.3.1, intake via this mechanism is probably minimal relative to other pathways. Also, COC absorbed by this pathway would have been accounted for in the tissue analyses from which it was concluded that risk is negligible.

6.3 Uncertainty Analysis

The assumption used for this assessment was that all site receptors were exposed to all site contaminants within an ecosystem, i.e., habitat was not delineated by source area. The following are uncertainties that may affect the outcome of risk characterization:

- Impacts to individual organisms on a site-wide basis are considered in this assessment as a predictor of impacts to populations. Generally, except for threatened and endangered species, assessments need only evaluate population effects. Evaluating risks to individual organisms tends to overestimate risks to populations.

- In estimating effects to red-tailed hawks from lead and copper. a conservative calculation was used, and no depuration losses were assumed.

- Applicability of literature-derived data depends on types of results presented and methods used to arrive at these results. Test endpoints produced by laboratory and field tests may be reported as formally defined toxicological endpoints or as less stringently defined measures of mortality or sublethal effect; variations in format introduce a source of error when subsumed into a single acceptable level value. Thus, seemingly equivalent values may in fact be significantly different due to differences in test protocols, test conditions, or responses of individual organisms.

- Regulatory standards, criteria, and/or toxicological data were not available for every COC (i.e., barium and manganese), and thus they could not be evaluated for potential impacts. These data gaps may cause an underestimate of risk because unevaluated COC could be unrecognized sources of risk.

- Tissue concentrations may only represent one animal or two or more animals of different species and may lead to either over- or underestimations of exposure to a population. No samples representative of avian tissue COC concentrations were sampled, and extrapolation from mammalian values were used. Also, no reference sources were sampled, and comparisons had to be made to literature-derived values.

- Synergistic, additive, and antagonistic interactions between chemicals were not explored.

6.4 Conclusions and Recommendations

The overall objective of this investigation was to conduct a BERA in accordance with federal and state guidance, which, when implemented, would allow for the selection of a cost-effective plan of action that ensured the protection of the public health and the environment. This

assessment determined the COC to ecological receptors and gave estimates as to the relative toxicity of the COC to both aquatic and terrestrial ecosystems present at the SS. This assessment sought to

- Describe the observed or potential magnitude of adverse ecological effect(s) at the SS and the primary cause(s) of the effect(s)
- Characterize the ecological consequences of the "no further action" remedial alternative.

The intrusive field activities conducted ensured that data gaps were identified and the necessary information collected to support these goals.

This BERA found localized areas with potentially elevated risks. These areas include the following:

- Station HMC-5 on Half Moon Creek may have a slightly impaired benthic community. This may be attributable to the chronic effects of aluminum in surface water and sediment COC or to natural conditions.
- The Beaver Pond, Half Moon Creek Overflow Pond, SA-2 (Clean Burn Area), SA-5 (HT-2 Well Area), and SA-6 (West Substation) had locally elevated concentrations of COC.

The assessment endpoints are meant to be protective of communities or populations. Therefore, the risk to individuals that spend all their time in the most highly contaminated areas is not considered in meeting the assessment endpoints. According to the toxicity test results, benchmarks, criteria, and calculations presented in this BERA, risks to the aquatic and terrestrial populations at the SS are negligible, and, therefore, the assessment endpoints and goals have been met. No action would be an acceptable alternative unless the desire is to be protective of individual organisms. It that case, removal of isolated areas of high contamination should be considered.

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Appendix A
Analytical Data

SALMON SITE

Appendix A:
Analytical Data

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Surface Water Quality February 1993

Parameter	Table A-1 Insitu Water Quality Parameters										
	BEP-1	BEP-1A	BEP-2	GAP-1	GRC-1	GRC-2	HHC-1	HHC-2	HMC-1	HMC-2	HMC-3
pH	5.99	5.03	4.50	5.56	3.90	5.12	4.70	5.56	5.15	5.54	4.30
Dissolved Oxygen (mg/L)	5.30	3.78	2.74	6.25	8.70	7.80	5.01	8.40	7.78	7.76	8.02
Oxidation Reduction Potential (mV)	109	106	131	135	107	125	-232	180	175	156	148
Conductivity (mS/cm)	0.057	0.046	0.043	0.142	0.015	0.015	0.033	0.015	0.020	0.027	0.02
Temperature (°C)	17.5	19.6	12.6	14.9	14.2	12.0	10.6	12.2	11.6	12.0	11.0
Turbidity (ntu)	5	16	10	0.70	-10	2	1	10	-10	-10	1
Depth (m)	0.50	0.50	0.25	0.10	0.76	0.76	0.076	0.11	0.38	0.91	0.53
Width (m)	--- [1]	--- [1]	--- [1]	NR*	4.6	6.1	0.91	1.1	1.4	5.5	5.5
Velocity (m/s)	0	0	0	0	0.37 [2]	0.24 [2]	0.15	0.55	0.12	0.18	0.37 [3]
Date	2/22/93	2/22/93	2/23/93	2/25/93	2/23/93	2/24/93	2/27/93	2/24/93	2/25/93	2/25/93	2/24/93

NR* = Not Recorded.

1 Pond dimensions are estimated at 260 m x 240 m.

2 At a depth of 0.76 m.

3 At a depth of 0.53 m.

Table A-1 Insitu Water Quality Parameters				
Parameter	HMC-4	HMC-5	HOP-1	REP-1
pH	3.94	5.45	4.53	6.32
Dissolved Oxygen (mg/L)	7.70	7.30	0.45	8.83
Oxidation Reduction Potential (mV)	112	-218	136	145
Conductivity (mS/cm)	0.017	0.027	0.505	0.025
Temperature (°C)	12.2	10.6	11.3	14.0
Turbidity (ntu)	-10	1	-10	-10
Depth (m)	0.74	0.91	2.0	NR*
Width (m)	5.5	6.7	NR*	NR*
Velocity (m/s)	0.67 [4]	0.37	0	0
Date	2/23/93	2/27/93	2/23/93	2/25/93

NR* = Not Recorded.

[4] At a depth of 0.74 m.

Compound (mg/Kg)	Table A-2.1 Metals								
	BEP-1-F-C-I	Q*	DV**	HMC-4-F-C-I	Q*	DV**	REP-1-F-C-1	Q*	DV**
Aluminum	2.3	B	J	5.3	B	J	13.4		---
Antimony	2.6	U	---	2.6	U	---	2.6	U	---
Arsenic	0.49	U	---	0.45	U	---	0.44	UW	UJ
Barium	6.1	B	J	66.2		---	2.2	B	J
Beryllium	0.04	U	---	0.05	B	U	0.04	U	---
Cadmium	0.13	U	---	0.13	U	---	0.13	U	---
Calcium	8020		---	34600		---	3690		---
Chromium	0.25	B	J	0.38	B	J	0.35	B	J
Cobalt	0.28	U	---	0.28	U	---	0.28	U	---
Copper	3.6		---	1.4		---	1.1	B	U
Iron	24.3		---	13.0		---	41.6		---
Lead	0.05	U	---	0.12	B	J	0.08	BW	J
Magnesium	351.0		---	681.0		---	291.0		---
Manganese	28.2		---	29.1		---	13.4		---
Mercury	0.15		U	0.43		---	0.04	U	---
Nickel	0.95	U	---	0.93	U	---	0.95	U	---
Potassium	3130		---	2790		---	2910		---
Selenium	0.25	UW	UJ	0.34	S	---	0.22	UW	UJ
Silver	0.17	U	---	0.16	U	---	0.17	U	---
Sodium	942.0		---	1250.0		---	750.0		---
Thallium	0.05	BW	J	0.05	UW	UJ	0.04	UW	UJ
Vanadium	0.17	U	---	0.17	U	---	0.17	U	---
Zinc	15.1		---	19.6		---	17.6		---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-2.2 Pesticides/PCB's								
	BEP-1-F-C	Q*	DV**	HMC-4-F-C-I	Q*	DV**	REP-1-F-C-I	Q*	DV**
Alpha-BHC	46	U	---	48	U	---	48	U	---
Beta-BHC	46	U	---	48	U	---	48	U	---
Delta-BHC	46	U	---	48	U	---	48	U	---
Gamma-BHC	46	U	---	48	U	---	48	U	---
Heptachlor	46	U	---	48	U	---	48	U	---
Aldrin	46	U	---	48	U	---	48	U	---
Heptachlor epoxide	46	U	---	48	U	---	48	U	---
Endosulfan I	46	U	---	48	U	---	48	U	---
Dieldrin	92	U	---	96	U	---	95	U	---
4,4'-DDE	92	U	---	96	U	---	95	U	---
Endrin	92	U	---	96	U	---	95	U	---
Endosulfan II	92	U	---	96	U	---	95	U	---
4,4'-DDD	92	U	---	96	U	---	95	U	---
Endosulfan sulfate	92	U	---	96	U	---	95	U	---
4,4'-DDT	92	U	---	96	U	---	95	U	---
Methoxychlor	460	U	---	480	U	---	480	U	---
Endrin ketone	92	U	---	96	U	---	95	U	---
Alpha-chlordane	460	U	---	480	U	---	480	U	---
Gamma-chlordane	460	U	---	480	U	---	480	U	---
Toxaphene	920	U	---	960	U	---	950	U	---
Aroclor-1016	460	U	---	480	U	---	480	U	---
Aroclor-1221	460	U	---	480	U	---	480	U	---
Aroclor-1232	460	U	---	480	U	---	480	U	---
Aroclor-1242	460	U	---	480	U	---	480	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-2.2 Pesticides/PCB's								
	BEP-1-F-C	Q*	DV**	HMC-4-F-C-I	Q*	DV**	REP-1-F-C-I	Q*	DV**
Aroclor-1248	460	U	---	480	U	---	480	U	---
Aroclor-1254	920	U	---	960	U	---	950	U	---
Aroclor-1260	920	U	---	960	U	---	950	U	---

Parameter (pCi/g)	Table A-2.3 Gross Alpha/Beta					
	BEP-1-F-C-I	DV**	HMC-4-F-C-I	DV**	REP-1-F-C-I	DV**
Alpha	0.05	---	0.20	---	ND	---
Sigma Error +/-	0.07	---	0.14	---	-	---
MDA*	0.12	U	0.21	U	0.16	U
Beta	0.57	---	0.94	---	1.29	---
Sigma Error +/-	0.08	---	0.11	---	0.13	---
MDA*	0.09	---	0.12	---	0.12	---

Table A-2.4 Gamma Spectroscopy												
Nuclide (pCi/g)	BEP-1-F-C-I	DV**	BEP-1A-F-C-I	DV**	BEP-2-F-C-I	DV**	GRC-1-F-C-I	DV**	HMC-4-F-C-I	DV**	REP-1-F-C-I	DV**
Cesium 137	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.30	U	0.10	U	0.20	U	0.10	U	0.10	U
Potassium 40	7.20	---	ND	---	ND	---	ND	---	ND	---	2.80	---
Sigma Error +/-	0.80	---	-	---	-	---	-	---	-	---	0.40	---
MDA*	7.20	---	5.50	U	4.90	U	3.50	U	3.10	U	0.24	---
Lead 212	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	0.40	U	0.20	U	0.30	U	0.20	U	0.20	U

Nuclide (pCi/g)	Table A-2.5 Tritium											
	BEP-1-F-C-I	DV**	BEP-1A-F-C-I	DV**	BEP-2-F-C-I	DV**	GRC-1-F-C-I	DV**	HMC-4-F-C-I	DV**	REP-1-F-C-I	DV**
Tritium	ND	---	0.49	---	0.12	---	0.55	---	0.24	---	ND	---
Sigma Error +/-	-	---	0.10	---	0.07	---	0.10	---	0.07	---	-	---
MDA*	0.08	U	0.16	---	0.14	U	0.17	---	0.14	---	0.11	U

Sample Number	Table A-3.1 Grain Size Analysis			
	% Moisture	% Sand (>0.0625 mm)	% Silt (0.0625 mm to 0.0039 mm)	% Clay (<0.0039 mm)
BEP-1-GS-G-I	66.5	11.1	65.7	23.2
BEP-1A-GS-G-I	25.0	40.8	34.5	24.7
BEP-2-GS-G-I	27.0	86.5	7.6	5.9
GAP-1-GS-G-I	79.1	42.6	33.9	23.5
GRC-1-GS-G-I	20.6	100.0	0.0	0.0
GRC-2-GS-G-I	23.4	100.0	0.0	0.0
HHC-1-GS-G-I	30.6	100.0	0.0	0.0
HHC-2-GS-G-I	18.7	96.3	1.8	1.9
HMC-1-GS-G-I	21.7	95.2	2.0	2.8
HMC-2-GS-G-I	29.3	93.2	3.8	3.0
HMC-3-GS-G-I	24.3	98.2	0.7	1.1
HMC-4-GS-G-I	21.6	100.0	0.0	0.0
HMC-4-GS-G-I-D	22.7	100.0	0.0	0.0
HMC-5-GS-G-I	33.6	96.3	1.8	1.9
HOP-1-GS-G-I	26.2	91.6	4.9	3.6
REP-1-GS-G-I	28.2	70.9	7.0	22.2

Table A-3.2 Metals

Compound (mg/Kg)	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I-D	Q*	DV**	GRC-2-SD-G-I	Q*	DV**
Aluminum	13000.0		---	4090.0		---	7650.0		---	6490.0		---	214.0		---	188.6		NP	84.4		---
Antimony	132.0	U	---	36.1	U	---	29.3	U	---	72.2	U	---	15.0	U	---	15.0	U	NP	14.5	U	---
Arsenic	7.7	B	J	3.0	B	J	3.6	B	J	1.5	B	J	0.3	U	---	0.0014	U	NP	0.3	U	---
Barium	181.0	B	J	27.3	B	J	21.1	B	J	125.0	B	J	5.9	B	J	4.1	B	NP	4.7	B	J
Beryllium	2.10	U	---	0.58	U	---	0.47	U	---	1.20	U	---	0.33	B	U	0.24	U	NP	0.23	U	---
Cadmium	9.00	B	J	1.80	U	---	1.40	U	---	3.50	U	---	0.73	U	---	0.73	U	NP	0.70	U	---
Calcium	5630	BE	J	1090	BE	J	885	BE	J	3620	BE	J	195	BE	J	305	B	NP	202	BE	J
Chromium	13.2	B	J	5.5	B	J	8.6		---	7.8	B	J	1.0	U	---	1.2	B	NP	0.9	U	---
Cobalt	14.3	U	---	3.9	U	---	3.2	U	---	13.2	B	U	1.6	U	---	1.6	U	NP	1.6	U	---
Copper	12.1	B	U	2.5	B	U	5.0	B	U	5.1	B	U	1.0	B	U	1.0	B	NP	1.0	B	U
Iron	13000	E	J	4700	E	J	6680	E	J	12300	E	J	303	E	J	215		NP	186	E	J
Lead	20.20		---	7.00		---	3.90		---	17.00		---	0.61	B	J	0.0034	B	NP	0.62	B	J
Magnesium	1380.0	B	J	279.0	B	J	363.0	B	J	518.0	B	J	58.8	B	J	52.5	B	NP	44.3	B	J
Manganese	583.0		J	214.0		J	28.2		J	136.0		J	85.1		J	49.8		NP	14.3		J
Mercury	0.60	U	---	0.16	U	---	0.13	U	---	0.32	U	---	0.07	U	---	0.07	U	NP	0.07	U	---
Nickel	47.6	U	---	13.0	U	---	10.6	U	---	26.0	U	---	5.4	U	---	5.4	U	NP	5.2	U	---
Potassium	6400	U	---	1750	U	---	1420	U	---	3490	U	---	727	U	---	727	U	NP	701	U	---
Selenium	2.40	UW	UJ	0.65	U	---	0.53	UW	UJ	1.30	U	---	0.27	U	---	0.0014	U	NP	0.26	U	---
Silver	8.30	U	---	2.30	U	---	1.80	U	---	4.50	U	---	0.95	U	---	0.95	U	NP	0.91	U	---
Sodium	1230.0	B	J	232.0	B	J	144.0	B	J	600.0	B	J	77.6	B		55.4	B	NP	66.4	B	J
Thallium	2.40	U	---	0.65	U	---	0.53	U	---	1.30	U	---	0.27	U	---	0.0014	U	NP	0.26	U	---
Vanadium	27.70	B	J	8.90	B	J	14.70	B	J	8.50	B	J	0.97	U	---	0.97	U	NP	0.94	U	---
Zinc	111.0		---	29.7		---	18.8		---	23.4	B	U	2.2	B	U	2.6	B	NP	2.3	B	U

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	GRC-2-SD-G-I-D	Q*	DV**	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**	HMC-4-SD-G-I	Q*	DV**
	Table A-3.2 Metals																				
Aluminum	93.7		NP	560.0		---	2180.0		---	195.0		---	649.0		---	211.0		---	1550.0		---
Antimony	14.5	U	NP	19.5	U	---	21.0	U	---	14.8	U	---	17.7	U	---	15.4	U	---	26.7	U	---
Arsenic	0.3	U	NP	0.4	UW	UJ	0.4	UW	UJ	0.3	UW	UJ	0.3	U	---	0.3	U	---	0.7	B	J
Barium	2.4	B	NP	16.0	B	J	34.3	B	J	2.8	B	U	12.7	B	J	7.3	B	J	51.7	B	J
Beryllium	0.23	U	NP	0.32	U	---	0.34	U	---	0.24	U	---	0.29	U	---	0.25	U	---	0.43	U	---
Cadmium	0.70	U	NP	0.95	U	---	1.00	U	---	0.72	U	---	1.70		---	0.75	U	---	1.30	U	---
Calcium	179	B	NP	451	BE	J	713	BE	J	190	BE	J	389	BE	J	183	BE	J	579	BE	J
Chromium	0.09	U	NP	2.2	B	J	2.4	B	J	0.9	U	---	1.5	B	J	1.0	U	---	2.6	B	J
Cobalt	1.6	U	NP	2.1	U	---	2.3	U	---	1.6	U	---	1.9	U	---	1.7	U	---	2.9	U	---
Copper	1.0	U	NP	1.4	B	U	1.4	U	---	1.0	U	---	1.2	U	---	1.1	B	U	2.5	B	U
Iron	125		NP	743	E	J	1930	E	J	102	E	J	319	E	J	269	E	J	1040	E	J
Lead	1.07		NP	0.90	B	J	2.40		---	1.10		---	2.00		---	2.20		---	5.00		---
Magnesium	39.8	B	NP	79.8	B	J	174.0	B	J	47.2	B	J	77.9	B	J	46.6	B	J	166.0	B	J
Manganese	6.3		NP	36.5		J	32.2		J	1.6	B	UJ	47.1		J	48.4		J	84.5		J
Mercury	0.06	U	NP	0.11	B	J	0.09	U	---	0.07	U	---	0.08	U	---	0.07	U	---	0.12	U	---
Nickel	5.2	U	NP	7.0	U	---	7.6	U	---	5.3	U	---	6.4	U	---	5.5	U	---	9.6	U	---
Potassium	701	U	NP	942	U	---	1020	U	---	714	U	---	857	U	---	743	U	---	1290	U	---
Selenium	0.26	U	NP	0.35	U	---	0.38	U	---	0.27	U	---	0.32	U	---	0.28	U	---	0.48	U	---
Silver	0.91	U	NP	1.20	U	---	1.30	U	---	0.93	U	---	1.10	U	---	0.97	U	---	1.70	U	---
Sodium	59.8	B	NP	78.0	B	J	96.8	B	J	70.3	B	J	73.2	B	J	60.7	B	U	128.0	B	J
Thallium	0.26	U	NP	0.35	U	---	0.38	U	---	0.27	U	---	0.32	U	---	0.28	U	---	0.48	U	---
Vanadium	0.94	U	NP	1.30	U	---	3.60	B	J	0.96	U	---	1.10	U	---	0.99	U	---	1.80	B	J
Zinc	1.6	B	NP	3.3	B	U	7.0	B	U	4.5	B	U	3.5	B	U	2.2	B	U	10.0		U

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	Table A-3.2 Metals								
	HMC-5-SD-G-I	Q*	DV**	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
Aluminum	836.0		---	1780.0		---	6480.0		---
Antimony	17.8	U	---	16.1	U	---	16.2	U	---
Arsenic	0.3	BW	J	0.9	B	J	2.1	B	J
Barium	34.0	B	J	11.0	B	J	19.9	B	J
Beryllium	0.29	U	---	0.29	B	U	0.26	U	---
Cadmium	0.87	U	---	0.78	U	---	0.79	U	---
Calcium	454	BE	J	417	BE	J	369	BE	J
Chromium	1.2	B	J	14.8		---	9.3		---
Cobalt	3.7	B	U	1.7	U	---	2.9	B	U
Copper	1.2	U	---	2.8	B	U	4.1	B	U
Iron	942	E	J	1280	E	J	19300	E	J
Lead	3.40		---	10.00		---	5.00		---
Magnesium	85.9	B	J	106.0	B	J	152.0	B	J
Manganese	90.8		J	17.5		J	45.8		J
Mercury	0.12	B	J	0.07	U	---	0.07	U	---
Nickel	6.4	U	---	5.8	U	---	5.8	U	---
Potassium	862	U	---	777	U	---	784	U	---
Selenium	0.32	U	---	0.29	UW	UJ	0.29	UW	UJ
Silver	1.10	U	---	1.00	U	---	1.00	U	---
Sodium	83.6	B	J	112.0	B	J	80.8	B	J
Thallium	0.32	U	---	0.29	U	---	0.29	U	---
Vanadium	1.40	B	J	3.20	B	J	43.00		---
Zinc	6.1	B	U	5.5	B	U	4.6	B	U

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.3 Pesticides/PCB's																							
	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-2-SD-G-I	Q*	DV**	GRC-2-SD-G-I-D	Q*	DV**
Alpha-BHC	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Beta-BHC	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Delta-BHC	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Gamma-BHC	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Heptachlor	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Aldrin	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Heptachlor epoxide	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Endosulfan I	190	U	NP	49	U	---	42	U	---	100	U	---	20	U	---	20	U	---	20	U	---
Dieldrin	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
4,4'-DDE	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
Endrin	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
Endosulfan II	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
4,4'-DDD	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
Endosulfan sulfate	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
4,4'-DDT	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
Methoxychlor	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---
Endrin ketone	380	U	NP	97	U	---	83	U	---	200	U	---	39	U	---	39	U	---	39	U	---
Alpha-chlordane	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---
Gamma-chlordane	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---
Toxaphene	3800	U	NP	970	U	---	830	U	---	2000	U	---	390	U	---	390	U	---	390	U	---
Aroclor-1016	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---
Aroclor-1221	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---
Aroclor-1232	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---
Aroclor-1242	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.3 Pesticides/PCB's																							
	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-2-SD-G-I	Q*	DV**	GRC-2-SD-G-I-D	Q*	DV**
Aroclor-1248	1900	U	NP	490	U	---	420	U	---	1000	U	---	200	U	---	200	U	---	200	U	---
Aroclor-1254	3800	U	NP	970	U	---	830	U	---	2000	U	---	390	U	---	390	U	---	390	U	---
Aroclor-1260	3800	U	NP	970	U	---	830	U	---	2000	U	---	390	U	---	390	U	---	390	U	---

Compound (ug/Kg)	Table A-3.3 Pesticides/PCB's																							
	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**	HMC-4-SD-G-I	Q*	DV**	HMC-5-SD-G-I	Q*	DV**
Alpha-BHC	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Beta-BHC	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Delta-BHC	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Gamma-BHC	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Heptachlor	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Aldrin	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Heptachlor epoxide	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Endosulfan I	22	U	---	27	U	---	20	U	---	25	U	---	21	U	---	38	U	---	25	U	---
Dieldrin	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
4,4'-DDE	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
Endrin	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
Endosulfan II	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
4,4'-DDD	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
Endosulfan sulfate	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
4,4'-DDT	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
Methoxychlor	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---
Endrin ketone	44	U	---	55	U	---	40	U	---	50	U	---	42	U	---	76	U	---	49	U	---
Alpha-chlordane	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---
Gamma-chlordane	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---
Toxaphene	440	U	---	550	U	---	400	U	---	500	U	---	420	U	---	760	U	---	490	U	---
Aroclor-1016	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---
Aroclor-1221	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---
Aroclor-1232	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---
Aroclor-1242	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---

Compound (ug/Kg)	Table A-3.3 Pesticides/PCB's																							
	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**	HMC-4-SD-G-I	Q*	DV**	HMC-5-SD-G-I	Q*	DV**
Aroclor-1248	220	U	---	270	U	---	200	U	---	250	U	---	210	U	---	380	U	---	250	U	---
Aroclor-1254	440	U	---	550	U	---	400	U	---	500	U	---	420	U	---	760	U	---	490	U	---
Aroclor-1260	440	U	---	550	U	---	400	U	---	500	U	---	420	U	---	760	U	---	490	U	---

Compound (ug/Kg)	Table A-3.3 Pesticides/PCB's					
	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
Alpha-BHC	22	U	---	23	U	---
Beta-BHC	22	U	---	23	U	---
Delta-BHC	22	U	---	23	U	---
Gamma-BHC	22	U	---	23	U	---
Heptachlor	22	U	---	23	U	---
Aldrin	22	U	---	23	U	---
Heptachlor epoxide	22	U	---	23	U	---
Endosulfan I	22	U	---	23	U	---
Dieldrin	44	U	---	46	U	---
4,4'-DDE	44	U	---	46	U	---
Endrin	44	U	---	46	U	---
Endosulfan II	44	U	---	46	U	---
4,4'-DDD	44	U	---	46	U	---
Endosulfan sulfate	44	U	---	46	U	---
4,4'-DDT	44	U	---	46	U	---
Methoxychlor	220	U	---	230	U	---
Endrin ketone	44	U	---	46	U	---
Alpha-chlordane	220	U	---	230	U	---
Gamma-chlordane	220	U	---	230	U	---
Toxaphene	440	U	---	460	U	---
Aroclor-1016	220	U	---	230	U	---
Aroclor-1221	220	U	---	230	U	---
Aroclor-1232	220	U	---	230	U	---
Aroclor-1242	220	U	---	230	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.3 Pesticides/PCB's					
	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
Aroclor-1248	220	U	---	230	U	---
Aroclor-1254	440	U	---	460	U	---
Aroclor-1260	440	U	---	460	U	---

Compound (ug/Kg)	Table A-3.4 Volatiles																							
	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-2-SD-G-I	Q*	DV**	GRC-2-SD-G-I-D	Q*	DV**
Chloromethane	120	U	---	32	U	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
Bromomethane	120	U	---	32	U	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
Vinyl Chloride	120	U	---	32	U	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
Chloroethane	120	U	---	32	U	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
Methylene Chloride	62	U	---	5	BJ	16U	13	U	---	33	U	---	7	U	---	2	J	---	2	J	---
Acetone	920	B		230	B		60	B	U	250	B	J	21	B	U	27	B	UJ	16	B	U
Carbon Disulfide	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
1,1-Dichloroethene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
1,1-Dichloroethane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
1,2-Dichloroethene (Total)	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Chloroform	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
1,2-Dichloroethane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
2-Butanone	170		---	5	J	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
1,1,1-Trichloroethane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Carbon Tetrachloride	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Vinyl Acetate	120	U	---	32	U	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
Bromodichloromethane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
1,2-Dichloropropane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
cis-1,3-Dichloropropene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Trichloroethene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Dibromochloromethane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
1,1,2-Trichloroethane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Benzene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---

* Data Qualifier. ** Data Validation.

Table A-3.4 Volatiles

Compound (ug/Kg)	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-2-SD-G-I	Q*	DV**	GRC-2-SD-G-I-D	Q*	DV**
trans-1,3-Dichloropropene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Bromoform	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
4-Methyl-2-Pentanone	120	U	---	32	U	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
2-Hexanone	120	U	---	32	U	---	26	U	---	67	U	---	14	U	---	13	U	---	13	U	---
Tetrachloroethene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
1,1,2,2-Tetrachlorethane	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Toluene	220		---	16	U	---	11	J	---	33	U	---	7	U	---	6	U	---	6	U	---
Chlorobenzene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Ethylbenzene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Styrene	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---
Xylene (Total)	62	U	---	16	U	---	13	U	---	33	U	---	7	U	---	6	U	---	6	U	---

Table A-3.4 Volatiles

Compound (ug/Kg)	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**	HMC-4-SD-G-I	Q*	DV**	HMC-5-SD-G-I	Q*	DV**
Chloromethane	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	---	16	U	---
Bromomethane	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	---	16	U	---
Vinyl Chloride	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	---	16	U	---
Chloroethane	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	---	16	U	---
Methylene Chloride	2	BJ	9U	3	J	---	1	J	---	8	U	UJ	7	U	---	8	BJ	12UJ	4	BJ	8U
Acetone	38	B	U	29	B	UJ	29	B	UJ	47	B	UJ	32	B	U	140	B	U	59	B	U
Carbon Disulfide	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
1,1-Dichloroethene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
1,1-Dichloroethane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
1,2-Dichloroethene (Total)	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
Chloroform	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
1,2-Dichloroethane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
2-Butanone	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	---	16	U	---
1,1,1-Trichloroethane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
Carbon Tetrachloride	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
Vinyl Acetate	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	---	16	U	---
Bromodichloromethane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
1,2-Dichloropropane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
cis-1,3-Dichloropropene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
Trichloroethene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
Dibromochloromethane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
1,1,2-Trichloroethane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
Benzene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---

* Data Qualifier. ** Data Validation.

Table A-3.4 Volatiles

Compound (ug/Kg)	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**	HMC-4-SD-G-I	Q*	DV**	HMC-5-SD-G-I	Q*	DV**
trans-1,3-Dichloropropene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
Bromoform	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	---	8	U	---
4-Methyl-2-Pentanone	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	UJ	16	U	---
2-Hexanone	18	U	---	19	U	---	13	U	---	16	U	UJ	14	U	---	24	U	UJ	16	U	---
Tetrachloroethene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	UJ	8	U	---
1,1,2,2-Tetrachlorethane	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	UJ	8	U	---
Toluene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	UJ	8	U	---
Chlorobenzene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	UJ	8	U	---
Ethylbenzene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	UJ	8	U	---
Styrene	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	UJ	8	U	---
Xylene (Total)	9	U	---	9	U	---	7	U	---	8	U	UJ	7	U	---	12	U	UJ	8	U	---

Compound (ug/Kg)	Table A-3.4 Volatiles					
	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
Chloromethane	14	U	---	14	U	---
Bromomethane	14	U	---	14	U	---
Vinyl Chloride	14	U	---	14	U	---
Chloroethane	14	U	---	14	U	---
Methylene Chloride	2	BJ	7U	2	J	---
Acetone	36	B	U	32	B	U
Carbon Disulfide	7	U	---	7	U	---
1,1-Dichloroethene	7	U	---	7	U	---
1,1-Dichloroethane	7	U	---	7	U	---
1,2-Dichloroethene (Total)	7	U	---	7	U	---
Chloroform	7	U	---	7	U	---
1,2-Dichloroethane	7	U	---	7	U	---
2-Butanone	14	U	---	14	U	---
1,1,1-Trichloroethane	7	U	UJ	7	U	---
Carbon Tetrachloride	7	U	UJ	7	U	---
Vinyl Acetate	14	U	UJ	14	U	---
Bromodichloromethane	7	U	UJ	7	U	---
1,2-Dichloropropane	7	U	UJ	7	U	---
cis-1,3-Dichloropropene	7	U	UJ	7	U	---
Trichloroethene	7	U	UJ	7	U	---
Dibromochloromethane	7	U	UJ	7	U	---
1,1,2-Trichloroethane	7	U	UJ	7	U	---
Benzene	7	U	UJ	7	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.4 Volatiles					
	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
trans-1,3-Dichloropropene	7	U	UJ	7	U	---
Bromoform	7	U	UJ	7	U	---
4-Methyl-2-Pentanone	14	U	---	14	U	---
2-Hexanone	14	U	---	14	U	---
Tetrachloroethene	7	U	---	7	U	---
1,1,2,2-Tetrachlorethane	7	U	---	7	U	---
Toluene	7	U	---	7	U	---
Chlorobenzene	7	U	---	7	U	---
Ethylbenzene	7	U	---	7	U	---
Styrene	7	U	---	7	U	---
Xylene (Total)	7	U	---	7	U	---

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles																	
	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-2-SD-G-I	Q*	DV**
Phenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
bis (2-Cloroethyl)Ether	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2-Chlorophenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
1,3-Dichlorobenzene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
1,4-Dichlorobenzene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Benzyl Alcohol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
1,2-Dichlorobenzene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2-Methylphenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
bis (2-Chloroisopropyl)Ether	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
4-Methylphenol	2200	J	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
N-Nitroso-Di-n-Propylamine	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Hexachloroethane	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Nitrobenzene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Isophorone	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2-Nitrophenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2,4-Dimethylphenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Benzoic Acid	20000	U	---	550	J	J	4200	U	---	10000	U	---	2000	U	---	2000	U	---
bis (2-Chloroethoxy)Methane	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2,4-Dichlorophenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
1,2,4-Trichlorobenzene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Naphthalene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
4-Chloroaniline	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Hexachlorobutadiene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
4-Chloro-3-Methylphenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles																	
	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-2-SD-G-I	Q*	DV**
2-Methylnaphthalene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Hexachlorocyclopentadiene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2,4,6-Trichlorophenol	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2,4,5-Trichlorophenol	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---
2-Chloronaphthalene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2-Nitroaniline	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---
Dimethyl Phthalate	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Acenaphthylene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2,6-Dinitrotoluene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
3-Nitroaniline	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---
Acenaphthene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2,4-Dinitrophenol	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---
4-Nitrophenol	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---
Dibenzofuran	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
2,4-Dinitrotoluene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Diethylphthalate	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
4-Chlorophenyl-phenylether	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Fluorene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
4-Nitroaniline	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---
4,6-Dinitro-2-Methylphenol	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---
N-Nitrosodiphenylamine (1)	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
4-Bromophenyl-phenylether	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Hexachlorobenzene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Pentachlorophenol	20000	U	---	4800	U	---	4200	U	---	10000	U	---	2000	U	---	2000	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles																	
	BEP-1-SD-G-I	Q*	DV**	BEP-1A-SD-G-I	Q*	DV**	BEP-2-SD-G-I	Q*	DV**	GAP-1-SD-G-I	Q*	DV**	GRC-1-SD-G-I	Q*	DV**	GRC-2-SD-G-I	Q*	DV**
Phenanthrene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Anthracene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Di-n-Butylphthalate	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Fluoranthene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Pyrene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Butylbenzylphthalate	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
3,3'-Dichlorbenzidine	8200	U	---	2000	U	---	1700	U	---	4100	U	---	810	U	---	820	U	---
Benzo(a)Anthracene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Chrysene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Bis (2-Ethylhexyl)Phthalate	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Di-n-Octyl Phthalate	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Benzo(b)Fluoranthene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Benzo(k)Fluoranthene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Benzo(a)Pyrene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Indeno(1,2,3-cd)Pyrene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Dibenz(a,h)Anthracene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---
Benzo(g,h,i)Perylene	4100	U	---	990	U	---	870	U	---	2100	U	---	410	U	---	410	U	---

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles																	
	GRC-2-SD-G-I-D	Q*	DV**	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**
Phenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
bis (2-Cloroethyl)Ether	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2-Chlorophenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
1,3-Dichlorobenzene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
1,4-Dichlorobenzene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Benzyl Alcohol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
1,2-Dichlorobenzene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2-Methylphenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
bis (2-Chloroisopropyl)Ether	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
4-Methylphenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
N-Nitroso-Di-n-Propylamine	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Hexachloroethane	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Nitrobenzene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Isophorone	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2-Nitrophenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2,4-Dimethylphenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Benzoic Acid	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
bis (2-Chloroethoxy)Methane	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2,4-Dichlorophenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
1,2,4-Trichlorobenzene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Naphthalene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
4-Chloroaniline	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Hexachlorobutadiene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
4-Chloro-3-Methylphenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles																	
	GRC-2-SD-G-I-D	Q*	DV**	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**
2-Methylnaphthalene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Hexachlorocyclopentadiene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2,4,6-Trichlorophenol	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2,4,5-Trichlorophenol	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
2-Chloronaphthalene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2-Nitroaniline	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
Dimethyl Phthalate	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Acenaphthylene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2,6-Dinitrotoluene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
3-Nitroaniline	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
Acenaphthene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2,4-Dinitrophenol	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
4-Nitrophenol	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
Dibenzofuran	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
2,4-Dinitrotoluene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Diethylphthalate	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
4-Chlorophenyl-phenylether	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Fluorene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
4-Nitroaniline	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
4,6-Dinitro-2-Methylphenol	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---
N-Nitrosodiphenylamine (1)	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
4-Bromophenyl-phenylether	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Hexachlorobenzene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Pentachlorophenol	2000	U	---	2200	U	---	2800	U	---	2000	U	---	2500	U	---	2100	U	---

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles																				
	GRC-2-SD-G-I-D	Q*	DV**	HHC-1-SD-G-I	Q*	DV**	HHC-2-SD-G-I	Q*	DV**	HMC-1-SD-G-I	Q*	DV**	HMC-2-SD-G-I	Q*	DV**	HMC-3-SD-G-I	Q*	DV**
Phenanthrene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Anthracene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Di-n-Butylphthalate	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Fluoranthene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Pyrene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Butylbenzylphthalate	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
3,3'-Dichlorbenzidine	820	U	---	900	U	---	1100	U	---	830	U	---	1000	U	---	850	U	---
Benzo(a)Anthracene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Chrysene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Bis (2-Ethylhexyl)Phthalate	410	U	---	450	U	---	570	U	---	420	U	---	.510	U	---	430	U	---
Di-n-Octyl Phthalate	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Benzo(b)Fluoranthene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Benzo(k)Fluoranthene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Benzo(a)Pyrene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Indeno(1,2,3-cd)Pyrene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Dibenz(a,h)Anthracene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---
Benzo(g,h,i)Perylene	410	U	---	450	U	---	570	U	---	420	U	---	510	U	---	430	U	---

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles											
	HMC-4-SD-G-I	Q*	DV**	HMC-5-SD-G-I	Q*	DV**	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
Phenol	780	U	---	510	U	---	460	U	---	480	U	---
bis (2-Cloroethyl)Ether	780	U	---	510	U	---	460	U	---	480	U	---
2-Chlorophenol	780	U	---	510	U	---	460	U	---	480	U	---
1,3-Dichlorobenzene	780	U	---	510	U	---	460	U	---	480	U	---
1,4-Dichlorobenzene	780	U	---	510	U	---	460	U	---	480	U	---
Benzyl Alcohol	780	U	---	510	U	---	460	U	---	480	U	---
1,2-Dichlorobenzene	780	U	---	510	U	---	460	U	---	480	U	---
2-Methylphenol	780	U	---	510	U	---	460	U	---	480	U	---
bis (2-Chloroisopropyl)Ether	780	U	---	510	U	---	460	U	---	480	U	---
4-Methylphenol	780	U	---	510	U	---	460	U	---	480	U	---
N-Nitroso-Di-n-Propylamine	780	U	---	510	U	---	460	U	---	480	U	---
Hexachloroethane	780	U	---	510	U	---	460	U	---	480	U	---
Nitrobenzene	780	U	---	510	U	---	460	U	---	480	U	---
Isophorone	780	U	---	510	U	---	460	U	---	480	U	---
2-Nitrophenol	780	U	---	510	U	---	460	U	---	480	U	---
2,4-Dimethylphenol	780	U	---	510	U	---	460	U	---	480	U	---
Benzoic Acid	3800	U	---	2500	U	---	2200	U	---	2300	U	---
bis (2-Chloroethoxy)Methane	780	U	---	510	U	---	460	U	---	480	U	---
2,4-Dichlorophenol	780	U	---	510	U	---	460	U	---	480	U	---
1,2,4-Trichlorobenzene	780	U	---	510	U	---	460	U	---	480	U	---
Naphthalene	780	U	---	510	U	---	460	U	---	480	U	---
4-Chloroaniline	780	U	---	510	U	---	460	U	---	480	U	---
Hexachlorobutadiene	780	U	---	510	U	---	460	U	---	480	U	---
4-Chloro-3-Methylphenol	780	U	---	510	U	---	460	U	---	480	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles											
	HMC-4-SD-G-I	Q*	DV**	HMC-5-SD-G-I	Q*	DV**	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
2-Methylnaphthalene	780	U	---	510	U	---	460	U	---	480	U	---
Hexachlorocyclopentadiene	780	U	---	510	U	---	460	U	---	480	U	---
2,4,6-Trichlorophenol	780	U	---	510	U	---	460	U	---	480	U	---
2,4,5-Trichlorophenol	3800	U	---	2500	U	---	2200	U	---	2300	U	---
2-Chloronaphthalene	780	U	---	510	U	---	460	U	---	480	U	---
2-Nitroaniline	3800	U	---	2500	U	---	2200	U	---	2300	U	---
Dimethyl Phthalate	780	U	---	510	U	---	460	U	---	480	U	---
Acenaphthylene	780	U	---	510	U	---	460	U	---	480	U	---
2,6-Dinitrotoluene	780	U	---	510	U	---	460	U	---	480	U	---
3-Nitroaniline	3800	U	---	2500	U	---	2200	U	---	2300	U	---
Acenaphthene	780	U	---	510	U	---	460	U	---	480	U	---
2,4-Dinitrophenol	3800	U	---	2500	U	---	2200	U	---	2300	U	---
4-Nitrophenol	3800	U	---	2500	U	---	2200	U	---	2300	U	---
Dibenzofuran	780	U	---	510	U	---	460	U	---	480	U	---
2,4-Dinitrotoluene	780	U	---	510	U	---	460	U	---	480	U	---
Diethylphthalate	780	U	---	510	U	---	460	U	---	480	U	---
4-Chlorophenyl-phenylether	780	U	---	510	U	---	460	U	---	480	U	---
Fluorene	780	U	---	510	U	---	460	U	---	480	U	---
4-Nitroaniline	3800	U	---	2500	U	---	2200	U	---	2300	U	---
4,6-Dinitro-2-Methylphenol	3800	U	---	2500	U	---	2200	U	---	2300	U	---
N-Nitrosodiphenylamine (1)	780	U	---	510	U	---	460	U	---	480	U	---
4-Bromophenyl-phenylether	780	U	---	510	U	---	460	U	---	480	U	---
Hexachlorobenzene	780	U	---	510	U	---	460	U	---	480	U	---
Pentachlorophenol	3800	U	---	2500	U	---	2200	U	---	2300	U	---

Compound (ug/Kg)	Table A-3.5 Semi-Volatiles											
	HMC-4-SD-G-I	Q*	DV**	HMC-5-SD-G-I	Q*	DV**	HOP-1-SD-G-I	Q*	DV**	REP-1-SD-G-I	Q*	DV**
Phenanthrene	780	U	---	510	U	---	460	U	---	480	U	---
Anthracene	780	U	---	510	U	---	460	U	---	480	U	---
Di-n-Butylphthalate	780	U	---	510	U	---	460	U	---	480	U	---
Fluoranthene	780	U	---	510	U	---	460	U	---	480	U	---
Pyrene	780	U	---	510	U	---	460	U	---	480	U	---
Butylbenzylphthalate	780	U	---	510	U	---	460	U	---	480	U	---
3,3'-Dichlorbenzidine	1600	U	---	1000	U	---	910	U	---	960	U	---
Benzo(a)Anthracene	780	U	---	510	U	---	460	U	---	480	U	---
Chrysene	780	U	---	510	U	---	460	U	---	480	U	---
Bis (2-Ethylhexyl)Phthalate	780	U	---	510	U	---	460	U	---	480	U	---
Di-n-Octyl Phthalate	780	U	---	510	U	---	460	U	---	480	U	---
Benzo(b)Fluoranthene	780	U	---	510	U	---	460	U	---	480	U	---
Benzo(k)Fluoranthene	780	U	---	510	U	---	460	U	---	480	U	---
Benzo(a)Pyrene	780	U	---	510	U	---	460	U	---	480	U	---
Indeno(1,2,3-cd)Pyrene	780	U	---	510	U	---	460	U	---	480	U	---
Dibenz(a,h)Anthracene	780	U	---	510	U	---	460	U	---	480	U	---
Benzo(g,h,i)Perylene	780	U	---	510	U	---	460	U	---	480	U	---

Table A-3.6 Gross Alpha/Beta																				
Parameter (pCi/g)	BEP-1-SD-G-I	DV**	BEP-1A-SD-G-I	DV**	BEP-2-SD-G-I	DV**	GAP-1-SD-G-I	DV**	GRC-1-SD-G-I	DV**	GRC-1-SD-G-I-D	DV**	GRC-2-SD-G-I	DV**	GRC-2-SD-G-I-D	DV**	HHC-1-SD-G-I	DV**
Alpha	ND	---	6.47	---	17.21	---	22.48	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	7.09	---	11.47	---	10.71	---	-	---	-	---	-	---	-	---	-	---
MDA*	15.21	U	11.42	U	17.10	---	12.30	J	11.15	U	15.05	U	13.40	UJ	10.70	UJ	13.32	U
Beta	18.57	---	15.20	---	11.13	---	23.69	---	11.29	---	8.35	---	9.29	---	11.66	---	5.90	---
Sigma Error +/-	5.76	---	4.69	---	5.05	---	6.00	---	4.25	---	8.46	---	5.29	---	4.20	---	4.18	---
MDA*	8.42	---	6.58	---	7.76	---	8.13	---	6.24	---	6.95	---	8.37	---	6.00	---	6.79	U

Parameter (pCi/g)	Table A-3.6 Gross Alpha/Beta																	
	HHC-2-SD-G-I	DV**	HMC-1-SD-G-I	DV**	HMC-2-SD-G-I	DV**	HMC-3-SD-G-I	DV**	HMC-4-SD-G-I	DV**	HMC-5-SD-G-I	DV**	HOP-1-SD-G-I	DV**	REP-1-SD-G-I	DV**
Alpha	ND	---	ND	---	10.20	---	5.88	---	7.48	---	6.63	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	7.69	---	7.18	---	7.18	---	5.51	---	-	---	-	---
MDA*	12.84	UJ	12.89	UJ	11.08	UJ	11.87	UJ	11.24	U	7.97	U	14.48	U	13.78	UJ
Beta	11.23	---	9.20	---	11.39	---	8.84	---	10.78	---	8.90	---	11.59	---	12.61	---
Sigma Error +/-	5.27	---	4.59	---	4.82	---	4.62	---	4.78	---	3.92	---	4.44	---	4.94	---
MDA*	8.10	---	7.17	---	7.35	---	7.29	---	7.36	---	5.99	---	6.48	---	7.63	---

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-3.7 Gamma Spectroscopy																					
	BEP-1-SD-G-I	DV**	BEP-1A-SD-G-I	DV**	BEP-2-SD-G-I	DV**	GAP-1-SD-G-I	DV**	GRC-1-SD-G-I	DV**	GRC-1-SD-G-I-D	DV**	GRC-2-SD-G-I	DV**	GRC-2-SD-G-I-D	DV**	HHC-1-SD-G-I	DV**				
Potassium 40	5.21	---	1.54	---	ND	---	2.38	---	0.41	---	0.25	---	ND	---	ND	---	-	---				
Sigma Error +/-	0.60	---	0.42	---	-	---	0.50	---	0.06	---	0.03	---	-	---	-	---	-	---				
MDA*	2.75	U	1.48	U	3.62	U	2.38	---	0.20	U	0.18	U	0.47	---	0.68	---	-	---				
Cesium 137	0.61	---	0.17	---	ND	---	1.09	---	0.013	---	0.012	---	ND	---	ND	---	ND	---				
Sigma Error +/-	0.05	---	0.02	---	-	---	0.08	---	0.003	---	0.002	---	-	---	-	---	-	---				
MDA*	0.10	---	0.06	---	0.14	U	0.08	---	0.009	---	0.008	---	0.02	---	0.03	---	0.20	---				
Lead 210	9.56	---	2.04	---	ND	---	-	---	0.24	---	0.31	---	0.24	---	-	---	-	---				
Sigma Error +/-	0.70	---	0.25	---	-	---	-	---	0.04	---	0.04	---	0.06	---	-	---	-	---				
MDA*	1.02	---	0.64	---	1.41	U	-	---	0.14	U	0.13	U	0.22	U	-	---	-	---				
Radium 224	ND	---	1.31	---	ND	---	4.14	---	0.24	---	ND	---	ND	---	ND	---	-	---				
Sigma Error +/-	-	---	0.39	---	-	---	0.58	---	0.06	---	-	---	-	---	-	---	-	---				
MDA*	2.30	---	0.87	---	2.01	U	1.82	---	0.15	U	0.13	U	0.28	---	0.49	---	-	---				
Radium 226	ND	---	0.68	---	0.36	---	ND	---	0.14	---	0.17	---	0.20	---	0.28	---	ND	---				
Sigma Error +/-	-	---	0.03	---	0.04	---	-	---	0.01	---	0.01	---	0.01	---	0.01	---	-	---				
MDA*	0.23	---	0.13	U	0.20	U	0.24	---	0.02	U	0.02	U	0.02	U	0.03	U	0.30	---				
Radium 228	ND	---	0.81	---	0.80	---	ND	---	0.15	---	0.14	---	0.17	---	0.22	---	-	---				
Sigma Error +/-	-	---	0.06	---	0.09	---	-	---	0.01	---	0.01	---	0.01	---	0.013	---	-	---				
MDA*	0.79	---	0.24	---	0.32	---	0.47	---	0.04	U	0.03	U	0.05	U	0.06	U	-	---				
Uranium 238	1.34	---	1.44	---	ND	---	ND	---	0.17	---	0.24	---	0.23	---	ND	---	-	---				
Sigma Error +/-	0.35	---	0.19	---	-	---	-	---	0.04	---	0.03	---	0.07	---	-	---	-	---				
MDA*	1.15	U	0.66	U	1.04	U	1.42	---	0.13	U	0.12	U	0.17	U	0.36	---	-	---				
Thallium 208	-	---	-	---	0.12	---	-	---	0.038	---	0.008	---	0.05	---	0.08	---	-	---				
Sigma Error +/-	-	---	-	---	0.03	---	-	---	0.004	---	-	---	0.01	---	0.01	---	-	---				
MDA*	-	---	-	---	0.10	---	-	---	0.01	---	-	---	0.01	U	0.02	U	-	---				

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-3.7 Gamma Spectroscopy																					
	BEP-1-SD-G-I	DV**	BEP-1A-SD-G-I	DV**	BEP-2-SD-G-I	DV**	GAP-1-SD-G-I	DV**	GRC-1-SD-G-I	DV**	GRC-1-SD-G-I-D	DV**	GRC-2-SD-G-I	DV**	GRC-2-SD-G-I-D	DV**	HHC-1-SD-G-I	DV**				
Lead 212	-	---	0.88	---	0.85	---	1.05	---	0.16	---	0.12	---	0.18	---	0.19	---	ND	---				
Sigma Error +/-	-	---	0.05	---	0.06	---	0.06	---	0.01	---	0.01	---	0.01	---	0.01	---	-	---				
MDA*	-	---	0.11	---	0.16	---	0.18	---	0.02	---	0.02	---	0.02	U	0.04	U	0.30	---				
Bismuth 212	-	---	-	---	-	---	-	---	0.11	---	-	---	-	---	-	---	-	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	0.03	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	0.07	---	-	---	-	---	-	---	-	---				
Cadmium 109	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	2.00	---				
Beryllium 7	2.40	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Sigma Error +/-	0.23	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	0.75	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Iodine 131	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Protactinium 231	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Protactinium 234	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	29.40	---				
Yttrium 88	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.20	---				

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-3.7 Gamma Spectroscopy																	
	BEP-1-SD-G-I	DV**	BEP-1A-SD-G-I	DV**	BEP-2-SD-G-I	DV**	GAP-1-SD-G-I	DV**	GRC-1-SD-G-I	DV**	GRC-1-SD-G-I-D	DV**	GRC-2-SD-G-I	DV**	GRC-2-SD-G-I-D	DV**	HHC-1-SD-G-I	DV**
Cerium 144	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.80	---
Thorium 230	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	18.90	---
Bismuth 214	0.43	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	0.06	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.27	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

Nuclide (pCi/g)	Table A-3.7 Gamma Spectroscopy																	
	HHC-2-SD-G-I	DV**	HMC-1-SD-G-I	DV**	HMC-2-SD-G-I	DV**	HMC-3-SD-G-I	DV**	HMC-4-SD-G-I	DV**	HMC-5-SD-G-I	DV**	HOP-1-SD-G-I	DV**	REP-1-SD-G-I	DV**
Potassium 40	0.94	---	0.57	---	0.26	---	0.39	---	0.90	---	-	---	1.42	---	1.50	---
Sigma Error +/-	0.12	---	0.07	---	0.03	---	0.05	---	0.09	---	-	---	0.30	---	0.18	---
MDA*	0.75	U	0.27	U	0.07	U	0.19	U	0.37	U	-	---	0.10	U	0.31	U
Cesium 137	0.07	---	ND	---	ND	---	ND	---	0.08	---	ND	---	0.16	---	ND	---
Sigma Error +/-	0.01	---	-	---	-	---	-	---	0.01	---	-	---	0.04	---	-	---
MDA*	0.02	---	0.21	---	0.01	---	0.01	---	0.02	---	0.10	---	0.03	---	0.03	---
Lead 210	-	---	0.41	---	-	---	0.26	---	1.15	---	-	---	0.88	---	0.99	---
Sigma Error +/-	-	---	0.06	---	-	---	0.05	---	0.10	---	-	---	0.08	---	0.10	---
MDA*	-	---	0.19	U	-	---	0.16	U	0.23	U	-	---	0.36	U	0.260	U
Radium 224	ND	---	0.30	---	ND	---	ND	---	0.68	---	-	---	ND	---	1.26	---
Sigma Error +/-	-	---	0.11	---	-	---	-	---	0.17	---	-	---	-	---	0.36	---
MDA*	0.61	---	0.20	U	0.17	---	0.22	---	0.22	---	-	---	0.77	U	0.24	---
Radium 226	0.46	---	0.31	---	ND	---	0.18	---	0.40	---	0.36	---	ND	---	0.76	---
Sigma Error +/-	0.02	---	0.01	---	-	---	0.01	---	0.02	---	0.03	---	-	---	0.03	---
MDA*	0.04	U	0.03	U	0.02	---	0.02	U	0.03	U	0.17	---	0.08	U	0.03	U
Radium 228	0.30	---	0.27	---	ND	---	0.15	---	0.37	---	-	---	ND	---	1.16	---
Sigma Error +/-	0.020	---	0.02	---	-	---	0.010	---	0.02	---	-	---	-	---	0.04	---
MDA*	0.07	U	0.05	U	0.03	---	0.03	U	0.05	U	-	---	0.16	U	0.04	---
Uranium 238	ND	---	0.36	---	0.27	---	0.17	---	0.33	---	-	---	0.67	---	1.21	---
Sigma Error +/-	-	---	0.05	---	0.03	---	0.06	---	0.06	---	-	---	0.14	---	0.11	---
MDA*	0.41	---	0.18	U	0.12	U	0.13	U	0.21	U	-	---	0.39	U	0.25	U
Thallium 208	0.13	---	-	---	-	---	0.043	---	-	---	-	---	0.11	---	0.39	---
Sigma Error +/-	0.01	---	-	---	-	---	0.004	---	-	---	-	---	0.01	---	0.02	---
MDA*	0.020	---	-	---	-	---	0.010	U	-	---	-	---	0.04	U	0.02	---

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-3.7 Gamma Spectroscopy																			
	HHC-2-SD-G-I	DV**	HMC-1-SD-G-I	DV**	HMC-2-SD-G-I	DV**	HMC-3-SD-G-I	DV**	HMC-4-SD-G-I	DV**	HMC-5-SD-G-I	DV**	HOP-1-SD-G-I	DV**	REP-1-SD-G-I	DV**				
Lead 212	0.39	---	ND	---	ND	---	0.19	---	-	---	0.43	---	0.34	---	1.42	---				
Sigma Error +/-	0.02	---	-	---	-	---	0.01	---	-	---	0.07	---	0.03	---	0.06	---				
MDA*	0.04	U	0.02	---	0.02	---	0.02	U	-	---	0.14	---	0.07	U	0.03	---				
Bismuth 212	-	---	-	---	-	---	0.09	---	-	---	-	---	-	---	-	---				
Sigma Error +/-	-	---	-	---	-	---	0.02	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	0.07	---	-	---	-	---	-	---	-	---				
Cadmium 109	-	---	0.74	---	-	---	-	---	0.67	---	ND	---	-	---	-	---				
Sigma Error +/-	-	---	0.14	---	-	---	-	---	0.13	---	-	---	-	---	-	---				
MDA*	-	---	0.20	---	-	---	-	---	0.29	U	1.10	---	-	---	-	---				
Beryllium 7	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
Iodine 131	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.10	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.04	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.01	---				
Protactinium 231	-	---	-	---	-	---	-	---	-	---	-	---	-	---	1.11	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.22	---				
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.47	---				
Protactinium 234	-	---	-	---	-	---	-	---	-	---	ND	---	-	---	3.89	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.66	---				
MDA*	-	---	-	---	-	---	-	---	-	---	20.50	---	-	---	2.96	---				
Yttrium 88	-	---	-	---	-	---	-	---	-	---	ND	---	-	---	-	---				
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---				
MDA*	-	---	-	---	-	---	-	---	-	---	0.12	---	-	---	-	---				

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-3.7 Gamma Spectroscopy																	
	HHC-2-SD-G-I	DV**	HMC-1-SD-G-I	DV**	HMC-2-SD-G-I	DV**	HMC-3-SD-G-I	DV**	HMC-4-SD-G-I	DV**	HMC-5-SD-G-I	DV**	HOP-1-SD-G-I	DV**	REP-1-SD-G-I	DV**
Cerium 144	-	---	-	---	-	---	-	---	-	---	0.30	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	0.10	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	0.30	---	-	---	-	---
Thorium 230	-	---	-	---	-	---	-	---	-	---	ND	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	11.10	---	-	---	-	---
Bismuth 214	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

Nuclide (pCi/g)	Table A-3.8 Tritium																			
	BEP-1-SD-G-I	DV**	BEP-1A-SD-G-I	DV**	BEP-2-SD-G-I	DV**	GAP-1-SD-G-I	DV**	GRC-1-SD-G-I	DV**	GRC-1-SD-G-I-D	DV**	GRC-2-SD-G-I	DV**	GRC-2-SD-G-I-D	DV**	HHC-1-SD-G-I	DV**
Tritium	0.48	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	0.10	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.17	---	0.16	U	0.11	U	0.17	U	0.04	U	0.04	U	0.06	U	0.06	U	0.06	U

Nuclide (pCi/g)	Table A-3.8 Tritium																	
	HHC-2-SD-G-I	DV**	HMC-1-SD-G-I	DV**	HMC-2-SD-G-I	DV**	HMC-3-SD-G-I	DV**	HMC-4-SD-G-I	DV**	HMC-5-SD-G-I	DV**	HOP-1-SD-G-I	DV**	REP-1-SD-G-I	DV**
Tritium	ND	---	0.03	---	ND	---	ND	---	ND	---	0.04	---	ND	---	ND	---
Sigma Error +/-	-	---	0.03	---	-	---	-	---	-	---	0.04	---	-	---	-	---
MDA*	0.06	U	0.06	U	0.08	U	0.06	U	0.13	U	0.08	U	0.10	U	0.07	U

* Minimum Detectable Activity. ** Data Validation.

Parameter (ug/g)	BEP-1-SD-G-I	DV**	BEP-1A-SD-G-I	DV**	BEP-2-SD-G-I	DV**	GAP-1-SD-G-I	DV**	GRC-1-SD-G-I	DV**	GRC-1-SD-G-I-D	DV**	GRC-2-SD-G-I	DV**	GRC-2-SD-G-I-D	DV**	HHC-1-SD-G-I	DV**
Table A-3.9 Ammonia and Total Organic Carbon																		
Ammonia	195	---	10.2	---	2.99	---	2.61	---	ND	---	ND	---	0.760	---	0.764	---	2.49	---
Detection Limit	17.7	---	0.971	---	0.789	---	1.94	---	0.404	---	0.404	---	0.389	---	0.389	---	0.521	---
Total Organic Carbon (TOC)	>190000	J	46700	J	27400	J	63800	J	674	J	585	J	365	J	397	J	18600	J
Detection Limit	295	---	80.9	---	65.7	---	161	---	33.6	---	33.6	---	32.4	---	32.4	---	43.6	---

** Data Validation.

Parameter (ug/g)	Table A-3.9 Wet Chemistry																	
	HHC-2-SD-G-I	DV**	HMC-1-SD-G-I	DV**	HMC-2-SD-G-I	DV**	HMC-3-SD-G-I	DV**	HMC-4-SD-G-I	DV**	HMC-5-SD-G-I	DV**	HOP-1-SD-G-I	DV**	REP-1-SD-G-I	DV**
Ammonia	1.40	---	ND	---	1.35	---	1.01	---	1.45	---	3.43	---	1.18	---	0.642	---
Detection Limit	0.565	---	0.397	---	0.476	---	0.412	---	0.718	---	0.478	---	0.431	---	0.436	---
Total Organic Carbon (TOC)	5870	J	4620	J	5520	J	2120	J	33300	J	11800	J	10700	J	9100	J
Detection Limit	47.1	---	33.1	---	39.6	---	34.4	---	59.9	---	39.9	---	35.9	---	36.3	---

** Data Validation.

Table A-4.1 Metals

Compound (ug/L)	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I-D	Q*	DV**	GRC-2-SW-G-I	Q*	DV**
Aluminum	261.0		---	228.0		---	416.0		---	110.0	B	J	179.0	B	J	164.8	B	NP	87.8	B	J
Antimony	55.6	U	---	55.6	U	---	55.6	U	---	55.6	U	---	55.6	U	---	55.6	U	NP	55.6	U	---
Arsenic	1.0	UW	UJ	1.0	UW	UJ	2.2	BW	J	1.0	U	---	1.0	UW	UJ	1.0	U	NP	1.0	U	---
Barium	31.3	B	J	26.2	B	J	18.9	B	J	32.6	B	J	25.6	B	J	25.1	B	NP	19.9	B	J
Beryllium	0.9	U	---	0.9	U	---	0.9	U	---	0.9	U	---	0.9	U	---	0.9	U	NP	0.9	U	---
Cadmium	2.7	U	---	2.7	U	---	2.7	U	---	2.7	U	---	2.7	U	---	3.8	B	NP	2.7	U	---
Calcium	2450	B	J	2460	B	J	3660	B	J	3630	B	J	3070	B	J	2975	B	NP	594	B	J
Chromium	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	NP	3.5	U	---
Cobalt	6.0	U	---	6.0	U	---	6.0	U	---	6.0	U	---	10.1	B	J	6.0	U	NP	6.0	U	---
Copper	3.7	U	---	3.7	U	---	3.7	U	---	3.7	U	---	3.7	U	---	3.7	U	NP	3.7	U	---
Iron	1440	E	J	1450	E	J	2690	E	J	252	E	J	654	E	J	591		NP	417	E	J
Lead	1.0	U	---	1.0	UW	UJ	1.2	B	J	1.0	U	---	1.2	B	J	1.1	B	NP	1.0	U	---
Magnesium	584	B	J	560	B	---	568	B	J	1040	B	J	812	B	J	785	B	NP	474	B	J
Manganese	155.0		---	120.0		---	150.0		---	84.2		---	116.0		---	121.5		NP	56.3		---
Mercury	0.1	U	---	0.1	U	---	0.1	U	---	0.2	B	J	0.1	U	---	0.1	U	NP	0.1	U	---
Nickel	20.0	U	---	20.0	U	---	20.0	U	---	20.0	U	---	20.0	U	---	20.0	U	NP	20.0	U	---
Potassium	2690	U	---	2690	U	---	2690	U	---	2690	U	---	2690	U	---	2690	U	NP	2690	U	---
Selenium	1.0	U	---	1.0	U	---	1.0	U	---	1.0	UW	UJ	1.0	U	---	1.0	U	NP	1.0	UW	UJ
Silver	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	NP	3.5	U	---
Sodium	2910	B	J	2700	B	J	2440	B	J	18200		---	2160	B	J	2094	B	NP	1740	B	J
Thallium	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	NP	1.0	U	---
Vanadium	3.6	U	---	3.6	U	---	3.6	U	---	3.6	U	---	3.6	U	---	3.6	U	NP	3.6	U	---
Zinc	4.8	B	U	7.0	B	U	6.8	B	U	3.8	B	U	13.1	B	U	13.8	B	NP	3.6	B	U

* Data Qualifier. ** Data Validation.

Table A-4.1 Metals

Compound (ug/L)	GRC-2-SW-G-I-D	Q*	DV**	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**
Aluminum	117.0	B	NP	201.0		---	165.0	B	J	190.0	B	J	151.0	B	J	113.0	B	J	141.0	B	J
Antimony	55.6	U	NP	55.6	U	---	55.6	U	---	55.6	U	---	55.6	U	---	55.6	U	---	55.6	U	---
Arsenic	1.0	U	NP	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---	1.0	UW	UJ
Barium	18.8	B	NP	90.4	B	J	29.0	B	J	29.8	B	J	27.9	B	J	25.4	B	J	22.3	B	J
Beryllium	0.9	U	NP	0.9	U	---	0.9	U	---	0.9	U	---	0.9	U	---	0.9	U	---	0.9	U	---
Cadmium	3.2	B	NP	2.7	U	---	2.7	U	---	2.7	U	---	2.7	U	---	2.7	U	---	2.7	U	---
Calcium	569	B	NP	827	B	J	615	B	J	573	B	J	747	B	J	694	B	J	634	B	J
Chromium	3.5	U	NP	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	---	3.5	U	---
Cobalt	6.0	U	NP	6.0	U	---	6.0	U	---	6.0	U	---	6.0	U	---	6.0	U	---	6.0	U	---
Copper	3.7	U	NP	3.7	U	---	3.7	U	---	3.7	U	---	3.7	U	---	3.7	U	---	3.7	U	---
Iron	424		NP	206	E	J	118	E	J	111	E	J	208	E	J	204	E	J	329	E	J
Lead	1.0	U	NP	1.0	UW	UJ	1.1	B	J	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---
Magnesium	464	B	NP	531	B	J	405	B	J	363	B	J	677	B	J	639	B	J	534	B	J
Manganese	55.8		NP	79.2		---	26.9		---	34.4		---	22.6		---	21.5		---	35.3		---
Mercury	0.1	U	NP	0.1	U	---	0.1	U	---	0.1	U	---	0.1	U	---	0.1	U	---	0.1	U	---
Nickel	20.0	U	NP	20.0	U	---	20.0	U	---	20.0	U	---	20.0	U	---	20.0	U	---	20.0	U	---
Potassium	2690	U	NP	2690	U	---	2690	U	---	2690	U	---	2690	U	---	2690	U	---	2690	U	---
Selenium	1.0	U	NP	1.0	UW	UJ	1.0	UW	UJ	1.0	UW	UJ	1.0	UW	UJ	1.0	UW	UJ	1.0	UW	UJ
Silver	3.5	U	NP	3.5	U	---	3.5	U	---	3.5	U	---	5.3	B	J	3.5	U	---	3.5	U	---
Sodium	1866	B	NP	1680	B	J	1640	B	J	1610	B	J	2110	B	J	2050	B	J	1950	B	J
Thallium	1.0	U	NP	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---	1.0	U	---
Vanadium	3.8	U	NP	3.6	U	---	3.6	U	---	3.6	U	---	3.6	U	---	3.6	U	---	3.6	U	---
Zinc	3.3	B	NP	5.8	B	U	5.1	B	J	4.3	B	U	5.2	B	J	4.3	B	U	3.0	B	U

* Data Qualifier. ** Data Validation.

Compound (ug/L)	Table A-4.1 Metals								
	HMC-5-SW-G-I	Q*	DV**	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-I	Q*	DV**
Aluminum	235.0		---	384.0		---	180.0	B	J
Antimony	55.6	U	---	55.6	U	---	55.6	U	---
Arsenic	1.0	U	---	1.0	UW	UJ	1.0	U	---
Barium	27.0	B	J	78.7	B	J	19.9	B	J
Beryllium	0.9	U	---	0.9	U	---	0.9	U	---
Cadmium	2.7	U	---	3.0	B	J	2.7	U	---
Calcium	798	B	J	24600		---	1620	B	J
Chromium	3.5	U	---	3.5	U	---	3.5	U	---
Cobalt	6.0	U	---	9.1	B	J	6.0	U	---
Copper	3.7	U	---	3.7	U	---	3.7	U	---
Iron	421	E	J	4190	E	J	677	E	J
Lead	1.0	UW	UJ	1.0	U	---	1.0	U	---
Magnesium	581	B	J	2220	B	J	566	B	J
Manganese	52.5		---	576.0		---	213.0		---
Mercury	0.1	U	---	0.1	U	---	0.2	B	J
Nickel	20.0	U	---	20.0	U	---	20.0	U	---
Potassium	2690	U	---	2690	U	---	2690	U	---
Selenium	1.0	UW	UJ	1.0	UW	UJ	1.0	U	---
Silver	3.5	B	J	3.5	U	---	14.6		---
Sodium	2120	B	J	56400		---	1220	B	J
Thallium	1.0	U	---	1.0	UW	UJ	1.0	U	---
Vanadium	3.6	U	---	3.6	U	---	3.6	U	---
Zinc	4.5	B	U	7.5	B	U	5.9	B	U

* Data Qualifier. ** Data Validation.

Compound (ug/L)	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-2-SW-G-I	Q*	DV**	GRC-2-SW-G-I-D	Q*	DV**
Table A-4.2 Pesticides/PCB's																					
Alpha-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Beta-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Delta-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Gamma-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Heptachlor	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Aldrin	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Heptachlor epoxide	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Endosulfan I	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Dieldrin	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
4,4'-DDE	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Endrin	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Endosulfan II	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
4,4'-DDD	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Endosulfan sulfate	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
4,4'-DDT	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Methoxychlor	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---
Endrin ketone	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Alpha-chlordane	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---
Gamma-chlordane	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---
Toxaphene	0.95	U	---	0.95	U	---	0.95	U	---	0.98	U	---	0.95	U	---	0.95	U	---	0.97	U	---
Aroclor-1016	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---
Aroclor-1221	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---
Aroclor-1232	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/L)	Table A-4.2 Pesticides/PCB's																							
	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-2-SW-G-I	Q*	DV**	GRC-2-SW-G-I-D	Q*	DV**			
Aroclor-1242	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---			
Aroclor-1248	0.48	U	---	0.48	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---			
Aroclor-1254	0.95	U	---	0.95	U	---	0.95	U	---	0.98	U	---	0.95	U	---	0.95	U	---	0.97	U	---			
Aroclor-1260	0.95	U	---	0.95	U	---	0.95	U	---	0.98	U	---	0.95	U	---	0.95	U	---	0.97	U	---			

Compound (ug/l)	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**	HMC-5-SW-G-I	Q*	DV**
	Table A-4.2 Pesticides/PCB's																				
Alpha-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Beta-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Delta-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Gamma-BHC	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Heptachlor	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Aldrin	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Heptachlor epoxide	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Endosulfan I	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Dieldrin	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
4,4'-DDE	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Endrin	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Endosulfan II	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
4,4'-DDD	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Endosulfan sulfate	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
4,4'-DDT	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Methoxychlor	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---
Endrin ketone	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---	0.10	U	---
Alpha-chlordane	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---
Gamma-chlordane	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---
Toxaphene	0.98	U	---	0.95	U	---	0.98	U	---	0.95	U	---	0.95	U	---	0.95	U	---	1.00	U	---
Aroclor-1016	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---
Aroclor-1221	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---
Aroclor-1232	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---

Compound (ug/L)	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**	HMC-5-SW-G-I	Q*	DV**
Table A-4.2 Pesticides/PCB's																					
Aroclor-1242	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---
Aroclor-1248	0.49	U	---	0.48	U	---	0.49	U	---	0.48	U	---	0.48	U	---	0.48	U	---	0.52	U	---
Aroclor-1254	0.98	U	---	0.95	U	---	0.98	U	---	0.95	U	---	0.95	U	---	0.95	U	---	1.00	U	---
Aroclor-1260	0.98	U	---	0.95	U	---	0.98	U	---	0.95	U	---	0.95	U	---	0.95	U	---	1.00	U	---

Compound (ug/L)	Table A-4.2 Pesticides/PCB's					
	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-I	Q*	DV**
Alpha-BHC	0.05	U	---	0.05	U	---
Beta-BHC	0.05	U	---	0.05	U	---
Delta-BHC	0.05	U	---	0.05	U	---
Gamma-BHC	0.05	U	---	0.05	U	---
Heptachlor	0.05	U	---	0.05	U	---
Aldrin	0.05	U	---	0.05	U	---
Heptachlor epoxide	0.05	U	---	0.05	U	---
Endosulfan I	0.05	U	---	0.05	U	---
Dieldrin	0.10	U	---	0.10	U	---
4,4'-DDE	0.10	U	---	0.10	U	---
Endrin	0.10	U	---	0.10	U	---
Endosulfan II	0.10	U	---	0.10	U	---
4,4'-DDD	0.10	U	---	0.10	U	---
Endosulfan sulfate	0.10	U	---	0.10	U	---
4,4'-DDT	0.10	U	---	0.10	U	---
Methoxychlor	0.48	U	---	0.48	U	---
Endrin ketone	0.10	U	---	0.10	U	---
Alpha-chlordane	0.48	U	---	0.48	U	---
Gamma-chlordane	0.48	U	---	0.48	U	---
Toxaphene	0.95	U	---	0.95	U	---
Aroclor-1016	0.48	U	---	0.48	U	---
Aroclor-1221	0.48	U	---	0.48	U	---
Aroclor-1232	0.48	U	---	0.48	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/L)	Table A-4.2 Pesticides/PCB's					
	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-I	Q*	DV**
Aroclor-1242	0.48	U	---	0.48	U	---
Aroclor-1248	0.48	U	---	0.48	U	---
Aroclor-1254	0.95	U	---	0.95	U	---
Aroclor-1260	0.95	U	---	0.95	U	---

Table A-4.3 Volatiles

Compound (ug/L)	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-2-SW-G-I	Q*	DV**	GRC-2-SW-G-I-D	Q*	DV**
Chloromethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bromomethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Vinyl Chloride	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	--	10	U	---
Chloroethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Methylene Chloride	2	BJ	5U	2	BJ	5U	2	BJ	5U	5	U	---	2	BJ	5U	5	U	---	5	U	---
Acetone	10	U	---	10	U	---	10	U	---	10	U	UJ	10	U	---	10	U	UJ	10	U	---
Carbon Disulfide	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethene (Total)	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Chloroform	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
2-Butanone	10	U	UJ	10	U	UJ	10	U	UJ	10	U	R	10	U	UJ	10	U	R	10	U	UJ
1,1,1-Trichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Carbon Tetrachloride	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Vinyl Acetate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bromodichloromethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloropropane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
cis-1,3-Dichloropropene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Trichloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Dibromochloromethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2-Trichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Benzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/L)	Table A-4.3 Volatiles																							
	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-2-SW-G-I	Q*	DV**	GRC-2-SW-G-I-D	Q*	DV**
trans-1,3-Dichloropropene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Bromoform	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
4-Methyl-2-Pentanone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Hexanone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Tetrachloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2,2-Tetrachlorethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Toluene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Chlorobenzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Ethylbenzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Styrene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Xylene (Total)	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---

Table A-4.3 Volatiles

Compound (ug/L)	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**	HMC-5-SW-G-I	Q*	DV**
Chloromethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bromomethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Vinyl Chloride	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Chloroethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Methylene Chloride	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	2	BJ	5U	5	U	---
Acetone	10	U	---	10	U	UJ	10	U	UJ	10	U	UJ	10	U	UJ	10	U	---	10	U	---
Carbon Disulfide	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethene (Total)	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Chloroform	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
2-Butanone	10	U	UJ	10	U	R	10	U	R	10	U	R	10	U	R	10	U	UJ	10	U	UJ
1,1,1-Trichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Carbon Tetrachloride	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Vinyl Acetate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bromodichloromethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloropropane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
cis-1,3-Dichloropropene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Trichloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Dibromochloromethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2-Trichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Benzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---

Compound (ug/L)	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**	HMC-5-SW-G-I	Q*	DV**
Table A-4.3 Volatiles																					
trans-1,3-Dichloropropene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Bromoform	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
4-Methyl-2-Pentanone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Hexanone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Tetrachloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2,2-Tetrachlorethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Toluene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Chlorobenzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Ethylbenzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Styrene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Xylene (Total)	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---

Compound (ug/L)	Table A-4.3 Volatiles																	
	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-1	Q*	DV**	Trip Blank 1	Q*	DV**	Trip Blank 2	Q*	DV**	Trip Blank 3	Q*	DV**	Trip Blank 4	Q*	DV**
Chloromethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bromomethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Vinyl Chloride	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Chloroethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Methylene Chloride	2	BJ	5U	5	U	---	2	BJ	5U	2	BJ	5U	2	U	---	5	U	---
Acetone	10	U	---	10	U	UJ	15		U	10	U	---	10	U	UJ	10	U	---
Carbon Disulfide	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethene (Total)	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Chloroform	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
2-Butanone	10	U	UJ	10	U	R	10	U	UJ	10	U	UJ	10	U	R	10	U	UJ
1,1,1-Trichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Carbon Tetrachloride	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Vinyl Acetate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bromodichloromethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloropropane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
cis-1,3-Dichloropropene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Trichloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Dibromochloromethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2-Trichloroethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Benzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---

Compound (ug/L)	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-1	Q*	DV**	Trip Blank 1	Q*	DV**	Trip Blank 2	Q*	DV**	Trip Blank 3	Q*	DV**	Trip Blank 4	Q*	DV**
	Table A-4.3 Volatiles																	
trans-1,3-Dichloropropene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Bromoform	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
4-Methyl-2-Pentanone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Hexanone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Tetrachloroethene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2,2-Tetrachlorethane	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Toluene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Chlorobenzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Ethylbenzene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Styrene	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---
Xylene (Total)	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---

Table A-4.4 Semi-Volatiles

Compound (ug/L)	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-2-SW-G-I	Q*	DV**	GRC-2-SW-G-I-D	Q*	DV**
Phenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
bis (2-Cloroethyl)Ether	10	U	UJ	10	U	UJ	10	U	UJ	10	U	---	10	U	UJ	10	U	---	10	U	---
2-Chlorophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,3-Dichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,4-Dichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzyl Alcohol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,2-Dichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Methylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
bis (2-Chloroisopropyl)Ether	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Methylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
N-Nitroso-Di-n-Propylamine	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Hexachloroethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Nitrobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Isophorone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Nitrophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dimethylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzoic Acid	48	U	---	48	U	---	48	U	---	50	U	UJ	48	U	---	48	U	UJ	48	U	UJ
bis (2-Chloroethoxy)Methane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dichlorophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,2,4-Trichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Naphthalene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Chloroaniline	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Hexachlorobutadiene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---

Compound (ug/L)	Table A-4.4 Semi-Volatiles																							
	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-2-SW-G-I	Q*	DV**	GRC-2-SW-G-I-D	Q*	DV**
4-Chloro-3-Methylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Methylnaphthalene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Hexachlorocyclopentadiene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4,6-Trichlorophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4,5-Trichlorophenol	48	U	---	48	U	---	48	U	---	50	U	---	48	U	---	48	U	---	48	U	---
2-Chloronaphthalene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Nitroaniline	48	U	---	48	U	---	48	U	---	50	U	---	48	U	---	48	U	---	48	U	---
Dimethyl Phthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Acenaphthylene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,6-Dinitrotoluene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
3-Nitroaniline	48	U	---	48	U	---	48	U	---	50	U	R	48	U	---	48	U	R	48	U	R
Acenaphthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dinitrophenol	48	U	---	48	U	---	48	U	---	50	U	---	48	U	---	48	U	---	48	U	---
4-Nitrophenol	48	U	---	48	U	---	48	U	---	50	U	---	48	U	---	48	U	---	48	U	---
Dibenzofuran	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dinitrotoluene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Diethylphthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Chlorophenyl-phenylether	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Fluorene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Nitroaniline	48	U	---	48	U	---	48	U	---	50	U	UJ	48	U	---	48	U	UJ	48	U	UJ
4,6-Dinitro-2-Methylphenol	48	U	---	48	U	---	48	U	---	50	U	---	48	U	---	48	U	---	48	U	---
N-Nitrosodiphenylamine (1)	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Bromophenyl-phenylether	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---

* Data Qualifier. ** Data Validation.

Table A-4.4 Semi-Volatiles

Compound (ug/L)	BEP-1-SW-G-I	Q*	DV**	BEP-1A-SW-G-I	Q*	DV**	BEP-2-SW-G-I	Q*	DV**	GAP-1-SW-G-I	Q*	DV**	GRC-1-SW-G-I	Q*	DV**	GRC-2-SW-G-I	Q*	DV**	GRC-2-SW-G-I-D	Q*	DV**
Hexachlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Pentachlorophenol	48	U	---	48	U	---	48	U	---	50	U	---	48	U	---	48	U	---	48	U	---
Phenanthrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Anthracene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Di-n-Butylphthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Fluoranthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Pyrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Butylbenzylphthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
3,3'-Dichlorbenzidine	19	U	---	19	U	---	19	U	---	20	U	---	19	U	---	19	U	---	19	U	---
Benzo(a)Anthracene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Chrysene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bis (2-Ethylhexyl)Phthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Di-n-Octyl Phthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(b)Fluoranthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(k)Fluoranthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(a)Pyrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Indeno(1,2,3-cd)Pyrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Dibenz(a,h)Anthracene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(g,h,i)Perylene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---

Compound (ug/L)	Table A-4.4 Semi-Volatiles																							
	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**	HMC-5-SW-G-I	Q*	DV**
Phenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
bis (2-Cloroethyl)Ether	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	UJ	10	U	---
2-Chlorophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,3-Dichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,4-Dichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzyl Alcohol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,2-Dichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Methylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
bis (2-Chloroisopropyl)Ether	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Methylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
N-Nitroso-Di-n-Propylamine	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Hexachloroethane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Nitrobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Isophorone	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Nitrophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dimethylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzoic Acid	48	U	UJ	48	U	UJ	49	U	UJ	48	U	UJ	48	U	UJ	48	U	---	49	U	UJ
bis (2-Chloroethoxy)Methane	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dichlorophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
1,2,4-Trichlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Naphthalene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Chloroaniline	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Hexachlorobutadiene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---

Table A-4.4 Semi-Volatiles

Compound (ug/L)	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**	HMC-5-SW-G-I	Q*	DV**
4-Chloro-3-Methylphenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Methylnaphthalene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Hexachlorocyclopentadiene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4,6-Trichlorophenol	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4,5-Trichlorophenol	48	U	---	48	U	---	49	U	---	48	U	---	48	U	---	48	U	---	49	U	---
2-Chloronaphthalene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2-Nitroaniline	48	U	---	48	U	---	49	U	---	48	U	---	48	U	---	48	U	---	49	U	---
Dimethyl Phthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Acenaphthylene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,6-Dinitrotoluene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
3-Nitroaniline	48	U	R	48	U	R	49	U	R	48	U	R	48	U	R	48	U	---	49	U	R
Acenaphthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dinitrophenol	48	U	---	48	U	---	49	U	---	48	U	---	48	U	---	48	U	---	49	U	---
4-Nitrophenol	48	U	---	48	U	---	49	U	---	48	U	---	48	U	---	48	U	---	49	U	---
Dibenzofuran	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
2,4-Dinitrotoluene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Diethylphthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Chlorophenyl-phenylether	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Fluorene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Nitroaniline	48	U	UJ	48	U	UJ	49	U	UJ	48	U	UJ	48	U	UJ	48	U	---	49	U	UJ
4,6-Dinitro-2-Methylphenol	48	U	---	48	U	---	49	U	---	48	U	---	48	U	---	48	U	---	49	U	---
N-Nitrosodiphenylamine (1)	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
4-Bromophenyl-phenylether	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---

Compound (ug/L)	Table A-4.4 Semi-Volatiles																							
	HHC-1-SW-G-I	Q*	DV**	HHC-2-SW-G-I	Q*	DV**	HMC-1-SW-G-I	Q*	DV**	HMC-2-SW-G-I	Q*	DV**	HMC-3-SW-G-I	Q*	DV**	HMC-4-SW-G-I	Q*	DV**	HMC-5-SW-G-I	Q*	DV**
Hexachlorobenzene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Pentachlorophenol	48	U	---	48	U	---	49	U	---	48	U	---	48	U	---	48	U	---	49	U	---	49	U	---
Phenanthrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Anthracene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Di-n-Butylphthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Fluoranthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Pyrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Butylbenzylphthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
3,3'-Dichlorbenzidine	19	U	UJ	19	U	---	19	U	---	19	U	---	19	U	---	19	U	---	20	U	UJ	20	U	UJ
Benzo(a)Anthracene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Chrysene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Bis (2-Ethylhexyl)Phthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Di-n-Octyl Phthalate	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(b)Fluoranthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(k)Fluoranthene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(a)Pyrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Indeno(1,2,3-cd)Pyrene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Dibenz(a,h)Anthracene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---
Benzo(g,h,i)Perylene	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---	10	U	---

Compound (ug/L)	Table A-4.4 Semi-Volatiles					
	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-I	Q*	DV**
Phenol	10	U	---	10	U	---
bis (2-Cloroethyl)Ether	10	U	UJ	10	U	---
2-Chlorophenol	10	U	---	10	U	---
1,3-Dichlorobenzene	10	U	---	10	U	---
1,4-Dichlorobenzene	10	U	---	10	U	---
Benzyl Alcohol	10	U	---	10	U	---
1,2-Dichlorobenzene	10	U	---	10	U	---
2-Methylphenol	10	U	---	10	U	---
bis (2-Chloroisopropyl)Ether	10	U	---	10	U	---
4-Methylphenol	10	U	---	10	U	---
N-Nitroso-Di-n-Propylamine	10	U	---	10	U	---
Hexachloroethane	10	U	---	10	U	---
Nitrobenzene	10	U	---	10	U	---
Isophorone	10	U	---	10	U	---
2-Nitrophenol	10	U	---	10	U	---
2,4-Dimethylphenol	10	U	---	10	U	---
Benzoic Acid	48	U	---	48	U	UJ
bis (2-Chloroethoxy)Methane	10	U	---	10	U	---
2,4-Dichlorophenol	10	U	---	10	U	---
1,2,4-Trichlorobenzene	10	U	---	10	U	---
Naphthalene	10	U	---	10	U	---
4-Chloroaniline	10	U	---	10	U	---
Hexachlorobutadiene	10	U	---	10	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/L)	Table A-4.4 Semi-Volatiles					
	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-I	Q*	DV**
4-Chloro-3-Methylphenol	10	U	---	10	U	---
2-Methylnaphthalene	10	U	---	10	U	---
Hexachlorocyclopentadiene	10	U	---	10	U	---
2,4,6-Trichlorophenol	10	U	---	10	U	---
2,4,5-Trichlorophenol	48	U	---	48	U	---
2-Chloronaphthalene	10	U	---	10	U	---
2-Nitroaniline	48	U	---	48	U	---
Dimethyl Phthalate	10	U	---	10	U	---
Acenaphthylene	10	U	---	10	U	---
2,6-Dinitrotoluene	10	U	---	10	U	---
3-Nitroaniline	48	U	---	48	U	R
Acenaphthene	10	U	---	10	U	---
2,4-Dinitrophenol	48	U	---	48	U	---
4-Nitrophenol	48	U	---	48	U	---
Dibenzofuran	10	U	---	10	U	---
2,4-Dinitrotoluene	10	U	---	10	U	---
Diethylphthalate	10	U	---	10	U	---
4-Chlorophenyl-phenylether	10	U	---	10	U	---
Fluorene	10	U	---	10	U	---
4-Nitroaniline	48	U	---	48	U	UJ
4,6-Dinitro-2-Methylphenol	48	U	---	48	U	---
N-Nitrosodiphenylamine (1)	10	U	---	10	U	---
4-Bromophenyl-phenylether	10	U	---	10	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/L)	Table A-4.4 Semi-Volatiles					
	HOP-1-SW-G-I	Q*	DV**	REP-1-SW-G-I	Q*	DV**
Hexachlorobenzene	10	U	---	10	U	---
Pentachlorophenol	48	U	---	48	U	---
Phenanthrene	10	U	---	10	U	---
Anthracene	10	U	---	10	U	---
Di-n-Butylphthalate	10	U	---	10	U	---
Fluoranthene	10	U	---	10	U	---
Pyrene	10	U	---	10	U	---
Butylbenzylphthalate	10	U	---	10	U	---
3,3'-Dichlorbenzidine	19	U	---	19	U	UJ
Benzo(a)Anthracene	10	U	---	10	U	---
Chrysene	10	U	---	10	U	---
Bis (2-Ethylhexyl)Phthalate	10	U	---	10	U	---
Di-n-Octyl Phthalate	10	U	---	10	U	---
Benzo(b)Fluoranthene	10	U	---	10	U	---
Benzo(k)Fluoranthene	10	U	---	10	U	---
Benzo(a)Pyrene	10	U	---	10	U	---
Indeno(1,2,3-cd)Pyrene	10	U	---	10	U	---
Dibenz(a,h)Anthracene	10	U	---	10	U	---
Benzo(g,h,i)Perylene	10	U	---	10	U	---

* Data Qualifier. ** Data Validation.

Table A-4.5 Gross Alpha/Beta																				
Parameter (pCi/L)	BEP-1-SW-G-I	DV**	BEP-1A-SW-G-I	DV**	BEP-2-SW-G-I	DV**	GAP-1-SW-G-I	DV**	GRC-1-SW-G-I	DV**	GRC-1-SW-G-I-D	DV**	GRC-2-SW-G-I	DV**	GRC-2-SW-G-I-D	DV**	HHC-1-SW-G-I	DV**	HHC-2-SW-G-I	DV**
Alpha	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	0.87	---	1.63	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.90	---	1.46	---	-	---
MDA*	2.50	UJ	3.51	UJ	2.61	UJ	4.44	UJ	2.41	UJ	2.46	UJ	2.51	UJ	1.35	UJ	2.17	UJ	2.32	UJ
Beta	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	3.58	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	3.59	---	-	---	-	---	-	---
MDA*	5.63	U	6.67	U	5.16	U	6.23	U	4.63	U	4.96	U	4.94	U	4.98	U	5.78	U	4.89	U

* Minimum Detectable Activity. ** Data Validation.

Parameter (pCi/L)	Table A-4.5 Gross Alpha/Beta															
	HMC-1-SW-G-I	DV**	HMC-2-SW-G-I	DV**	HMC-3-SW-G-I	DV**	HMC-4-SW-G-I	DV**	HMC-5-SW-G-I	DV**	HOP-1-SW-G-I	DV**	REP-1-SW-G-I	DV**
Alpha	1.16	---	ND	---	1.58	---	ND	---	ND	---	3.50	---	ND	---
Sigma Error +/-	1.14	---	-	---	1.29	---	-	---	-	---	3.21	---	-	---
MDA*	1.75	UJ	2.36	UJ	1.89	UJ	3.26	UJ	2.23	UJ	4.72	UJ	2.90	UJ
Beta	ND	---	ND	---	ND	---	ND	---	ND	---	2.51	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	3.27	---	-	---
MDA*	4.54	U	5.19	U	4.64	U	5.18	U	4.36	U	5.52	U	5.55	U

Nuclide (pCi/L)	Table A-4.6 Gamma Spectroscopy																			
	BEP-1-SW-G-I	DV**	BEP-1A-SW-G-I	DV**	BEP-2-SW-G-I	DV**	GAP-1-SW-G-I	DV**	GRC-1-SW-G-I	DV**	GRC-1-SW-G-1-D	DV**	GRC-2-SW-G-I	DV**	GRC-2-SW-G-I-D	DV**	HHC-1-SW-G-I	DV**
Tin 113	-	---	-	---	-	---	ND	---	-	---	-	---	8.70	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	2.10	---	-	---	-	---
MDA*	-	---	-	---	-	---	13.80	U	-	---	-	---		---	15.40	U	11.00	U
Cesium137	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	13.05	U	10.97	U	9.14	U	11.40	U	14.81	U	13.43	U	8.20	U	13.60	U	11.00	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/L)	Table A-4.6 Gamma Spectroscopy																	
	HHC-2-SW-G-I	DV**	HMC-1-SW-G-I	DV**	HMC-2-SW-G-I	DV**	HMC-3-SW-G-I	DV**	HMC-4-SW-G-I	DV**	HMC-5-SW-G-I	DV**	HOP-1-SW-G-I	DV**	REP-1-SW-G-I	DV**
Tin 113	ND	---	ND	---	ND	---	ND	---	-	---	ND	---	-	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	8.50	U	7.80	U	11.90	U	7.10	U	-	---	10.20	U	-	---	9.30	U
Cesium137	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	5.90	U	6.90	U	10.30	U	7.30	U	13.79	U	10.40	U	10.00	U	7.50	U

Table A-4.7 Tritium																				
Parameter (pCi/mL)	BEP-1-SW-G-I	DV**	BEP-1A-SW-G-I	DV**	BEP-2-SW-G-I	DV**	GAP-1-SW-G-I	DV**	GRC-1-SW-G-I	DV**	GRC-1-SW-G-I-D	DV**	GRC-2-SW-G-I	DV**	GRC-2-SW-G-I-D	DV**	HHC-1-SW-G-I	DV**
Tritium	0.40	---	0.40	---	0.20	---	ND	---	0.20	---	0.20	---	0.20	---	0.10	---	ND	---
Sigma Error +/-	0.12	---	0.11	---	0.10	---	-	---	0.11	---	0.11	---	0.10	---	0.09	---	-	---
MDA*	0.22	---	0.22	---	0.23	U	0.19	U	0.20	---	0.20	---	0.19	---	0.19	U	0.20	U

Table A-4.7 Tritium

Parameter (pCi/mL)	HHC-2-SW-G-I	DV**	HMC-1-SW-G-I	DV**	HMC-2-SW-G-I	DV**	HMC-3-SW-G-I	DV**	HMC-4-SW-G-I	DV**	HMC-5-SW-G-I	DV**	HOP-1-SW-G-I	DV**	REP-1-SW-G-I	DV**
Tritium	ND	---	0.10	---	0.10	---	ND	---	0.20	---	ND	---	0.70	---	ND	---
Sigma Error +/-	-	---	0.13	---	0.15	---	-	---	0.10	---	-	---	0.15	---	-	---
MDA*	0.19	U	0.19	U	0.19	U	0.19	U	0.20	---	0.20	U	0.23	---	0.19	U

Table A-4.8 Water Parameters																				
Category	BEP-1-SW-G-I	DV**	BEP-1A-SW-G-I	DV**	BEP-2-SW-G-I	DV**	GAP-1-SW-G-I	DV**	GRC-1-SW-G-I	DV**	GRC-1-SW-G-I-D	DV**	GRC-2-SW-G-I	DV**	GRC-2-SW-G-I-D	DV**	HHC-1-SW-G-I	DV**
Total Alkalinity (mg/L)	ND	---	ND	---	5.43	---	ND	---	ND	---	-	NP	ND	---	ND	---	ND	---
Ammonia (µg/L)	ND	---	ND	---	ND	---	33.20	---	ND	---	-	NP	35.30	---	ND	---	ND	---
Hardness (mg/L)	13.90	---	13.50	---	16.70	---	19.90	---	6.77	---	6.57	---	8.27	---	8.27	---	8.82	U
Nitrate/Nitrite (µg/L)	ND	---	ND	---	ND	---	ND	---	42.20	---	-	NP	36.00	---	33.50	---	ND	---
Ortho-Phosphate (µg/L)	45.00	---	36.80	---	39.80	---	69.70	---	59.20	---	-	NP	73.00	---	81.60	---	83.10	---
Total Phosphorus (µg/L)	ND	---	ND	---	ND	---	54.30	---	35.50	---	-	NP	ND	---	ND	---	35.50	---
TDS (mg/L)	35.00	---	35.00	---	38.00	---	67.00	---	6.00	---	-	NP	25.00	---	29.00	---	ND	---
TSS (mg/L)	4.00	---	5.00	---	8.00	---	260.00	---	6.00	---	-	NP	13.00	---	11.00	---	22.00	---

** Data Validation.

Table A-4.8 Wet Chemistry																						
Category	HHC-2-SW-G-I	DV**	HMC-1-SW-G-I	DV**	HMC-2-SW-G-I	DV**	HMC-3-SW-G-I	DV**	HMC-4-SW-G-I	DV**	HMC-5-SW-G-I	DV**	HOP-1-SW-G-I	DV**	REP-1-SW-G-I	DV**						
Total Alkalinity (mg/L)	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	5.33	---	ND	---						
Ammonia (µg/L)	ND	---	ND	---	ND	---	ND	---	34.70	---	ND	---	94.40	---	76.90	---						
Hardness (mg/L)	7.88	---	9.65	---	10.20	---	10.20	---	7.76	---	11.80	U	78.40	---	13.60	---						
Nitrate/Nitrite (µg/L)	ND	---	216.00	---	ND	---	198.00	---	113.00	---	112.00	---	ND	---	21.10	---						
Ortho-Phosphate (µg/L)	83.90	---	93.60	---	72.70	---	65.20	---	60.00	---	63.30	---	47.30	---	42.10	---						
Total Phosphorus (µg/L)	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---						
TDS (mg/L)	43.00	---	14.00	---	34.00	---	23.00	---	22.00	---	36.00	---	310.00	---	18.00	---						
TSS (mg/L)	9.00	---	3.00	---	6.00	---	3.00	---	4.00	---	10.00	---	18.00	---	7.00	---						

** Data Validation.

Parameter (mg/L)	Table A-4.9 Biochemical Oxygen Demand																	
	BEP-1-SW-G-I	DV**	BEP-1A-SW-G-I	DV**	GAP-1-SW-G-I	DV**	GRC-1-SW-G-I	DV**	GRC-1-SWBOD	DV**	GRC-2-SW-G-I	DV**	GRC-2-SW-G-I-D	DV**	HHC-1-SW-G-I	DV**	HHC-2-SWBOD	DV**
BOD	5.58	---	5.58	---	7.38	---	5.04	---	2.78	---	6.00	---	6.00	---	4.08	---	3.24	---

** Data Validation.

Parameter (mg/L)	Table A-4.9 Biochemical Oxygen Demand																	
	HMC-1-SW-G-I	DV**	HMC-2-SW-G-I	DV**	HMC-2-SWBOD	DV**	HMC-3-SW-G-I	DV**	HMC-4-SW-G-I	DV**	HMC-4-SWBOD	DV**	HMC-5-SW-G-I	DV**	HOP-1-SW-G-I	DV**	HOP-1-SWBOD	DV**
BOD	4.74	---	5.70	---	4.26	---	21.20	---	5.58	---	3.18	---	6.48	---	6.18	---	3.72	---

** Data Validation.

Table A-4.9 Biochemical Oxygen Demand

Parameter (mg/L)	REP-1-SW-G-I	DV**
BOD	8.28	---

** Data Validation.

Surface Water Sampling 1993

Table A-4.10

Summary of 48-Hour Ceriodaphnid Acute Toxicity Tests

Conducted April 27 - 29, 1993

Sample ID	Survival (%) Replicates A	B	C	D	E	Average	Significant [1] Difference vs. Reference
CONTROL (laboratory)	80	100	100	100	100	96	NA
Half Moon Creek Stations							
HMC-1 (reference)	80	100	100	100	60	88	NA
HMC-2	100	100	100	100	100	100	---
HMC-3	40	100	100	100	100	88	---
HMC-4	80	100	100	100	80	92	---
HMC-5	100	80	100	100	80	92	---
Grantham Creek Stations							
GRC-1 (reference)	100	40	100	100	100	88	NA
GRC-2	100	60	80	80	80	80	---
Hickory Hollow Creek Stations							
HHC-1 (reference)	80	40	60	60	60	60	NA
HHC-2	100	40	80	60	40	64	---
Pond Stations							
REP-1 (reference)	80	60	100	80	80	80	NA
BEP-1	40	80	100	80	60	72	---
BEP-1A	40	60	60	80	60	60	---
BEP-2	40	60	40	100	60	60	---
GAP-1	80	60	80	40	40	60	---
HOP-1	40	60	40	80	80	60	---

[1] An asterisk (*) indicates sample survival is significantly less than respective reference survival at the 95% confidence level.

A dash (-) indicates no difference.

Surface Water Sampling 1993

Table A-4.11

Summary of 48-Hour Fathead Minnow Acute Toxicity Tests

Conducted April 27 - 29, 1993

Sample ID	Survival (%) Replicates A	B	C	D	Average	Significant [1] Difference vs. Reference
CONTROL (laboratory)	100	100	100	100	100	NA
Half Moon Creek Stations						
HMC-1 (reference)	100	100	100	100	100	NA
HMC-2	100	100	100	100	100	---
HMC-3	100	80	100	100	95	---
HMC-4	100	100	100	100	100	---
HMC-5	100	100	100	100	100	---
Grantham Creek Stations						
GRC-1 (reference)	100	100	80	100	95	NA
GRC-2	100	100	100	100	100	---
Hickory Hollow Creek Stations						
HHC-1 (reference)	100	100	100	100	100	NA
HHC-2	100	100	100	100	100	---
Pond Stations						
REP-1 (reference)	100	100	100	100	100	NA
BEP-1	100	100	100	100	100	---
BEP-1A	100	100	100	100	100	---
BEP-2	100	100	100	100	100	---
GAP-1	100	100	100	100	100	---
HOP-1	100	100	100	100	100	---

[1] An asterisk (*) indicates sample survival is significantly less than respective reference survival at the 95% confidence level.

A dash (-) indicates no difference.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA1-1A-SS	Q*	DV**	SA1-1A-SS-D	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1D-SS-D	Q*	DV**	SA1-1E-SS	Q*	DV**
Aluminum	3480		J	2570		NP	5310		J	2450		J	2540		---	3690		NP	2470		J
Antimony	13.4	UN	UJ	13.4	U	NP	12.6	UN	UJ	12.4	UN	UJ	12.8	UN	UJ	12.8	U	NP	12.3	UN	UJ
Arsenic	1.90	BW	J	1.90	B	NP	1.60	B	J	1.50	B	J	1.60	B	J	1.60	B	NP	1.40	B	J
Barium	28.4	B	J	31.7	B	NP	14.9	B	J	31.4	B	J	11.9	B	J	15.2	B	NP	17.4	B	J
Beryllium	0.07	U	---	0.14	B	NP	0.06	U	---	0.13	B	U	0.07	B	U	0.14	B	NP	0.06	U	---
Cadmium	0.71	U	---	0.71	U	NP	0.67	U	---	0.66	U	---	0.68	U	---	0.68	U	NP	0.65	U	---
Calcium	1570		---	1380		NP	615	B	J	539	B	J	247	B	J	310	B	NP	798	B	J
Chromium	16.7		---	15.8		NP	44.0		---	4.0		---	4.2		---	5.3		NP	7.2		---
Cobalt	2.80	B	J	1.00	U	NP	2.20	B	J	1.20	B	J	1.10	B	J	1.40	B	NP	1.40	B	J
Copper	5.10	B	U	3.55	B	NP	2.90	B	U	3.30	B	U	0.63	B	U	2.75	B	NP	8.20		---
Iron	6540		J	4380		NP	11400		J	4950		J	5300		J	6100		NP	5370		J
Lead	16.6	N	J	13.3		NP	9.0	SN	J	14.5	N	J	4.3	N	J	4.0		NP	4.7	N	J
Magnesium	185.0	B	J	224.0	B	NP	155.0	B	J	103.0	B	J	82.0	B	J	116.0	B	NP	115.0	B	J
Manganese	48.8		J	30.7		NP	20.7		J	14.4		J	16.4		J	31.5		NP	22.4		J
Mercury	0.06	U	---	0.06	U	NP	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	NP	0.05	U	---
Nickel	4.8	U	---	4.8	U	NP	4.5	U	---	4.4	U	---	4.6	U	---	4.6	U	NP	13.1		---
Potassium	679	U	---	679	U	NP	641	U	---	858	B	J	878	B	J	650	U	NP	835	B	J
Selenium	0.35	B	U	0.23	U	NP	0.34	B	U	0.21	U	---	0.22	U	---	0.22	U	NP	0.21	U	---
Silver	0.80	U	---	0.80	U	NP	0.75	U	---	0.74	U	---	0.76	U	---	0.76	U	NP	0.73	U	---
Sodium	78.0	B	J	46.8	B	NP	53.7	B	J	31.3	B	U	47.7	B	U	51.6	B	NP	58.0	B	J
Thallium	0.23	U	---	0.23	U	NP	0.21	UW	UJ	0.21	U	---	0.22	U	---	0.22	U	NP	0.21	U	---
Vanadium	13.4		---	10.3	B	NP	25.1		---	10.3	B	J	13.7		---	14.8		NP	12.4		---
Zinc	15.0		---	13.9		NP	7.0		---	20.3		---	2.1	B	U	4.4	B	NP	6.5		---
Cyanide	0.15	U	---	0.15	U	NP	0.37	B	J	0.14	U	---	0.14	U	---	0.14	U	NP	0.14	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**
Aluminum	2540		J	3520		J	2200		J	2690		J	4860		J	2690		J	5090		J
Antimony	12.9	UN	UJ	13.6	UN	UJ	12.7	UN	UJ	13.8	UN	UJ	13.5	UN	---	15.0	UN	UJ	12.6	UN	UJ
Arsenic	1.80	B	J	2.00	B	J	2.10	B	J	3.20		---	11.90		---	1.80	B	J	1.80	B	J
Barium	6.1	B	J	7.8	B	J	7.7	B	J	14.0	B	J	16.9	B	J	19.8	B	J	10.7	B	J
Beryllium	0.07	U	---	0.07	U	---	0.06	U	---	0.07	U	---	0.14	B	U	0.08	U	---	0.13	B	U
Cadmium	0.68	U	---	0.72	U	---	0.67	U	---	0.73	U	---	0.71	U	---	0.79	U	---	0.67	U	---
Calcium	162	B	J	172	B	J	230	B	J	249	B	J	667	B	J	184	B	J	182	B	J
Chromium	6.9		---	7.2		---	5.4		---	5.8		---	16.6		---	4.6		---	9.6		---
Cobalt	1.30	B	J	1.90	B	J	1.60	B	J	2.20	B	J	3.00	B	J	1.60	B	J	1.70	B	J
Copper	1.30	B	U	5.40	B	U	0.85	B	U	1.60	B	---	1.40	B	U	1.00	B	U	2.10	B	U
Iron	6400		J	6610		J	5130		J	5260		J	6960		J	4530		J	9080		J
Lead	2.8	N	J	5.2	N	J	3.5	N	J	16.7	N	J	5.9	N	UJ	9.7	N	J	5.2	N	UJ
Magnesium	83.1	B	J	97.2	B	J	106.0	B	J	109.0	B	J	190.0	B	J	97.9	B	J	138.0	B	J
Manganese	9.1		J	13.2		J	14.9		J	23.1		J	47.3		J	10.5		J	26.5		J
Mercury	0.05	U	---	0.06	U	---	0.05	U	---	0.06	U	---	0.29		---	0.06	U	---	0.05	U	---
Nickel	4.6	U	---	4.8	U	---	4.5	U	---	4.9	U	---	4.8	U	---	5.3	U	---	4.5	U	---
Potassium	654	U	---	689	U	---	644	U	---	702	U	---	687	U	---	761	U	---	641	U	---
Selenium	0.22	U	---	0.23	U	---	0.22	U	---	0.24	U	---	0.23	U	---	0.26	UW	UJ	0.21	U	---
Silver	0.77	U	---	0.81	U	---	0.76	U	---	0.82	U	---	0.81	U	---	0.89	U	---	0.75	U	---
Sodium	37.4	B	U	33.2	B	U	45.4	B	U	41.1	B	U	49.5	B	U	45.7	B	U	30.4	B	U
Thallium	0.22	U	---	0.23	U	---	0.22	U	---	0.24	U	---	0.23	U	---	0.26	U	---	0.21	U	---
Vanadium	14.6		---	14.7		---	11.6		---	11.4	B	U	13.6		---	11.1	B	J	17.9		---
Zinc	2.5	B	U	4.2	B	U	5.3		---	6.0		---	23.0		---	5.3		---	7.5		---
Cyanide	0.35	B	J	0.15	U	---	0.14	U	---	0.15	U	---	0.15	U	---	0.16	U	---	0.14	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**	SA1-3E-SS	Q*	DV**	SA1-4A-SS	Q*	DV**	SA1-4A-SS-D	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**
Aluminum	2580		J	3880		J	1120		J	7060		---	6210		NP	4280		---	7510		---
Antimony	12.4	UN	UJ	13.5	UN	UJ	12.2	UN	UJ	14.5	UN	J	14.5	U	NP	13.6	UN	J	12.7	UN	J
Arsenic	0.54	B	U	2.30		---	1.30	B	J	1.10	BW	J	1.70	B	NP	1.50	BW	J	1.90	BW	J
Barium	11.1	B	J	18.9	B	J	4.5	B	J	76.9		---	68.0		NP	152.0		---	23.5	B	J
Beryllium	0.06	U	---	0.14	B	U	0.06	U	---	0.15	B	U	0.29	B	NP	0.07	B	U	0.07	B	U
Cadmium	0.65	U	---	0.71	U	---	0.65	U	---	0.77	U	---	0.77	U	NP	0.72	U	---	0.67	U	---
Calcium	259	B	J	344	B	J	144	B	J	439	B	J	427	B	NP	2230		---	323	B	J
Chromium	6.9		---	15.0		---	2.9		---	7.6		---	8.0		NP	6.3		---	7.1		---
Cobalt	1.30	B	J	1.90	B	J	1.40	B	J	3.80	B	U	5.20	B	NP	2.60	B	U	2.70	B	U
Copper	8.80		---	2.90	B	U	0.82	B	U	3.50	B	---	3.30	B	NP	5.10	B	J	5.40	B	J
Iron	3670		J	5600		J	3530		J	4560		---	4560		NP	5280		---	7000		---
Lead	18.5	N	UJ	9.6	N	J	7.8	SN	J	10.5		J	10.8		NP	11.0		J	4.5		J
Magnesium	116.0	B	J	167.0	B	J	49.8	B	J	183.0	B	J	151.0	B	NP	162.0	B	J	128.0	B	J
Manganese	21.3		J	66.3		J	6.5		J	636.0		---	600.0		NP	85.2		---	28.5		---
Mercury	0.05	U	---	0.06	U	---	0.05	U	---	0.06	U	---	0.06	U	NP	0.06	U	---	0.05	U	---
Nickel	4.4	U	---	4.8	U	---	4.3	U	---	6.5	B	J	5.3	B	NP	4.8	U	---	4.5	U	---
Potassium	629	U	---	685	U	---	621	U	---	738	U	---	738	U	NP	692	U	---	644	U	---
Selenium	0.21	U	---	0.23	U	---	0.21	U	---	0.25	UW	J	0.44	B	NP	0.23	U	---	0.22	U	---
Silver	0.74	U	---	0.80	U	---	0.73	U	---	0.87	U	---	0.87	U	NP	0.81	U	---	0.76	U	---
Sodium	46.7	B	U	52.3	B	U	38.2	B	U	48.3	B	U	49.9	B	NP	71.6	B	U	63.9	B	U
Thallium	0.21	U	---	0.23	U	---	0.21	U	---	0.25	U	---	0.25	U	NP	0.23	U	---	0.22	U	---
Vanadium	7.8	B	J	12.9		---	8.0	B	J	10.6	B	J	10.6	B	NP	10.5	B	J	15.1		---
Zinc	8.0		---	16.3		---	2.8	B	U	9.1		---	8.2		NP	28.8		---	10.3		---
Cyanide	0.14	U	---	0.15	U	---	0.25	B	J	0.15	U	---	0.15	U	NP	0.17	B	J	0.23	B	J

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**	SA2-5A-SS-D	Q*	DV**	SA2-5B-SS	Q*	DV**	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**
Aluminum	4710		---	6900		---	4670		---	4440		NP	4190		---	9210		---	3430		---
Antimony	13.7	UN	J	14.6	UN	J	13.8	UN	UJ	13.8	U	NP	13.6	UN	UJ	14.0	UN	UJ	12.9	UN	UJ
Arsenic	1.20	BW	J	2.60	S	---	1.30	BW	J	1.40	B	NP	2.20	BW	J	4.00	S	---	1.40	BW	J
Barium	27.2	B	J	208.0		---	43.8	B	J	42.4	B	NP	38.4	B	J	22.4	B	J	19.6	B	J
Beryllium	0.07	U	---	0.07	B	J	0.27	B	J	0.27	B	NP	0.27	B	J	0.28	B	J	0.07	U	---
Cadmium	0.72	U	---	0.77	U	---	0.73	U	---	0.73	U	NP	0.72	U	---	0.74	U	---	0.68	U	---
Calcium	250	B	J	1350		---	814	BE	J	785	B	NP	1790	E	J	676	BE	J	406	BE	J
Chromium	6.9		---	13.8		---	6.9		---	5.0		NP	6.6		---	7.7		---	3.9		---
Cobalt	3.60	B	U	5.90	B	J	7.10	B	J	8.80	B	NP	4.10	B	U	3.70	B	U	3.50	B	U
Copper	2.70	B	U	8.80		---	0.61	UN	---	0.61	U	NP	1.20	BN	J	0.62	UN	---	0.57	UN	---
Iron	4640		---	6110		---	4750		---	4730		NP	7810		---	11500		---	4270		---
Lead	5.9		J	16.0		J	10.2		---	10.9		NP	18.4		---	10.0		---	7.3		---
Magnesium	122.0	B	J	156.0	B	J	252.0	B	J	228.0	B	NP	351.0	B	J	615.0	B	J	183.0	B	J
Manganese	247.0		---	581.0		---	956.0		---	994.0		NP	482.0		---	31.1		---	274.0		---
Mercury	0.08	B	U	0.06	U	---	0.11	B	J	0.12		NP	0.07	B	J	0.06	U	---	0.06	U	---
Nickel	5.8	B	---	5.2	U	---	4.9	U	---	4.9	U	NP	4.8	U	---	5.0	U	---	4.6	U	---
Potassium	696	U	---	765	B	J	699	U	---	699	U	NP	690	U	---	710	U	---	657	U	---
Selenium	0.23	U	---	0.25	U	---	0.26	B	U	0.23	U	NP	0.23	U	---	0.24	U	---	0.22	U	---
Silver	0.82	U	---	0.87	U	---	0.82	U	---	0.82	U	NP	0.81	U	---	0.83	U	---	0.77	U	---
Sodium	40.0	B	U	154.0	B	J	71.2	B	U	53.5	B	NP	46.9	B	U	70.3	B	U	70.5	B	U
Thallium	0.23	U	---	0.25	U	---	0.23	U	---	0.23	U	NP	0.23	U	---	0.27	B	J	0.22	UW	UJ
Vanadium	9.9	B	J	11.4	B	J	9.2	B	J	9.4	B	NP	11.0	B	J	17.3		---	8.5	B	J
Zinc	8.5		---	31.8		---	7.3	N	J	19.5		NP	11.9	N	J	9.1	N	UJ	8.5	N	J
Cyanide	0.13	B	---	0.14	U	---	0.15	U	---	0.15	U	NP	0.18	B	J	0.15	U	---	0.14	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**	SA2-6D-SS	Q*	DV**	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**
Aluminum	6360		---	3730		---	2950		---	1610		---	2890		---	1910		---	5050		---
Antimony	13.6	UN	UJ	13.6	UN	UJ	12.6	UN	R	12.2	UN	R	12.7	UN	R	12.9	UN	R	13.7	U	---
Arsenic	2.20	BW	J	0.75	B	J	0.82	B	J	1.60	B	J	0.78	B	J	2.20	B	J	2.50		---
Barium	26.3	B	J	26.6	B	J	7.5	B	J	4.2	B	J	6.1	B	J	14.2	B	J	41.6	B	J
Beryllium	0.27	B	J	0.07	U	---	0.06	U	---	0.06	U	---	0.07	U	---	0.07	U	---	0.37	B	U
Cadmium	0.72	U	---	0.72	U	---	1.40		---	0.64	U	---	0.67	U	---	0.68	U	---	0.72	U	---
Calcium	409	BE	J	484	B	J	299	B	J	204	B	J	150	B	J	255	B	J	572	B	J
Chromium	6.9		---	6.4		---	4.5		---	3.8		---	4.7		---	2.2	B	J	12.4		---
Cobalt	9.30	B	J	6.00	B	J	1.10	B	J	2.30	B	J	1.60	B	J	2.30	B	J	7.20	B	J
Copper	0.60	UN	---	1.40	B	J	1.50	B	U	1.30	B	U	1.80	B	U	0.57	U	---	4.70	BE	UJ
Iron	9260		---	5130		---	4890		---	3720		---	5270		---	3940		---	6700		---
Lead	12.1		---	8.1		---	4.6		---	18.4		---	12.3		---	8.5		---	14.5		---
Magnesium	286.0	B	J	190.0	B	J	94.9	B	J	68.7	B	J	83.4	B	J	84.2	B	J	248.0	B	J
Manganese	202.0		---	564.0		---	41.6		---	21.1		---	18.7		---	278.0		---	696.0		J
Mercury	0.06	U	---	0.06	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.06	U	---
Nickel	4.8	U	---	4.8	U	---	4.5	U	---	4.3	U	---	4.5	U	---	4.6	U	---	4.9	U	---
Potassium	691	U	---	690	U	---	641	U	---	620	U	---	646	U	---	653	U	---	697	U	---
Selenium	0.23	U	---	0.56	BN	UJ	0.21	UN	UJ	0.21	UN	UJ	0.22	UN	UJ	0.23	BN	UJ	0.23	UWN	UJ
Silver	0.81	U	---	0.81	U	---	0.75	U	---	0.73	U	---	0.76	U	---	0.77	U	---	0.82	U	---
Sodium	44.1	B	U	49.0	B	U	44.0	B	U	38.7	B	U	47.4	B	U	25.8	B	U	44.1	B	U
Thallium	0.26	BW	J	0.23	U	---	0.21	U	---	0.21	U	---	0.22	U	---	0.22	U	---	0.23	U	---
Vanadium	17.4		---	10.6	B	J	10.5	B	J	8.0	B	J	10.6	B	J	8.4	B	J	14.1		---
Zinc	5.9	N	UJ	4.2	B	J	5.2		---	4.1	B	J	3.8	B	U	2.7	B	U	5.3		U
Cyanide	0.15	U	---	0.15	U	---	0.14	U	---	0.14	B	J	0.14	U	---	0.14	U	---	0.13	U	---

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**	SA3-8A-SS	Q*	DV**	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**
	Table A-5.1 Metals/Cyanide																				
Aluminum	4830		---	2080		---	2950		---	2450		---	5470		---	2770		---	1300		---
Antimony	14.2	U	---	12.8	U	---	13.0	U	---	12.4	U	---	14.2	U	---	12.5	U	---	12.1	U	---
Arsenic	2.80		---	1.50	B	J	4.60		---	6.10		---	2.50		---	1.50	B	J	0.49	B	J
Barium	42.4	B	J	12.7	B	J	29.2	B	J	12.7	B	J	74.0		---	23.2	B	J	7.0	B	J
Beryllium	0.30	B	U	0.07	B	U	0.08	B	U	0.07	B	J	0.67	B	U	0.06	U	---	0.07	B	U
Cadmium	0.75	U	---	0.67	U	---	0.68	U	---	0.66	U	---	0.75	U	---	0.66	U	---	0.64	U	---
Calcium	544	B	J	232	B	J	1230		---	215	B	J	686	B	J	3370		---	247	B	J
Chromium	11.5		---	5.3		---	16.4		---	5.9		---	7.7		---	5.1		---	2.1		---
Cobalt	8.50	B	J	2.30	B	U	5.00	B	U	3.30	B	U	10.60	B	J	3.10	B	U	1.40	B	U
Copper	20.50	E	UJ	12.00	E	UJ	9.70	E	UJ	5.70	E	UJ	5.60	BE	UJ	16.40	E	UJ	3.50	BE	UJ
Iron	10100		---	4660		---	15400		---	4480		---	5300		---	7340		---	2740		---
Lead	26.6		---	24.4	W	J	85.2	S	---	40.3		---	20.1		---	12.5		---	5.6		---
Magnesium	261.0	B	J	102.0	B	J	219.0	B	J	110.0	B	J	365.0	B	J	174.0	B	J	77.9	B	J
Manganese	746.0		J	109.0		J	865.0		J	183.0		J	977.0		J	118.0		J	33.7		J
Mercury	0.06	U	---	0.05	U	---	0.06	U	---	0.05	U	---	0.07	B	J	0.05	U	---	0.05	U	---
Nickel	5.0	U	---	4.5	U	---	5.9	B	J	4.4	U	---	5.0	U	---	4.4	U	---	4.3	U	---
Potassium	720	U	---	648	U	---	658	U	---	631	U	---	720	U	---	633	U	---	615	U	---
Selenium	0.24	UN	UJ	0.22	UN	UJ	0.22	UN	UJ	0.21	UWN	UJ	0.24	UN	UJ	0.21	UWN	UJ	0.21	UN	UJ
Silver	0.84	U	---	0.76	U	---	0.77	U	---	0.74	U	---	0.84	U	---	0.74	U	---	0.72	U	---
Sodium	61.5	B	U	43.9	B	U	47.5	B	U	51.1	B	U	61.5	B	U	49.4	B	U	41.7	B	U
Thallium	0.24	UW	UJ	0.22	U	---	0.22	U	---	0.21	U	---	0.24	UW	UJ	0.21	U	---	0.21	U	---
Vanadium	16.6		---	8.5	B	J	13.0		---	9.7	B	J	11.2	B	J	10.4	B	J	6.1	B	J
Zinc	54.3		---	17.5		---	278.0		---	76.4		---	8.0		---	12.2		---	3.5	B	U
Cyanide	0.14	U	---	0.13	U	---	0.15	U	---	0.13	U	---	0.16	U	---	0.13	U	---	0.12	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**	SA3-9C-SS	Q*	DV**	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**
Aluminum	2880		---	9560		---	18800		---	9010		---	3250		---	6180		---	8180		---
Antimony	12.3	U	---	13.8	U	---	14.7	U	---	13.9	U	---	12.8	U	---	14.6	U	---	14.4	U	---
Arsenic	1.80	B	J	3.20		---	3.20		---	2.30	B	J	2.40		---	1.90	B	J	2.30	B	J
Barium	31.4	B	J	48.9		---	72.7		---	50.9		---	84.1		---	60.3		---	236.0		---
Beryllium	0.07	B	U	0.44	B	U	0.86	B	U	0.52	B	U	0.21	B	U	0.39	B	U	0.39	B	U
Cadmium	0.65	U	---	0.73	U	---	0.77	U	---	0.74	U	---	0.83	B	U	0.77	U	---	0.76	U	---
Calcium	377	B	J	801	B	J	667	B	J	1000	B	J	1870		---	1710		---	2550		---
Chromium	10.8		---	10.8		---	18.2		---	13.9		---	7.5		---	6.7		---	13.5		---
Cobalt	5.90	B	U	7.30	B	J	7.10	B	U	6.80	B	U	1.70	B	U	7.70	B	J	7.80	B	J
Copper	11.60	E	UJ	6.10	E	UJ	13.30	E	UJ	9.00	E	UJ	5.70	E	UJ	10.40	E	UJ	8.40	E	UJ
Iron	10300		---	12900		---	16200		---	13900		---	6930		---	7530		---	10400		---
Lead	18.7	W	J	12.5	S	---	13.9	S	---	9.6	S	---	13.9		---	15.7		---	24.2		---
Magnesium	163.0	B	J	647.0	B	---	1180.0	B	J	517.0	B	J	211.0	B	J	337.0	B	J	460.0	B	J
Manganese	212.0		---	614.0		J	39.4		J	390.0		J	90.0		J	945.0		J	631.0		J
Mercury	0.05	U	---	0.06	U	---	0.06	U	---	0.07	B	J	0.06	B	J	0.06	U	---	0.06	B	J
Nickel	4.4	U	---	4.9	U	---	11.8		---	5.0	U	---	4.6	U	---	5.2	U	---	5.1	U	---
Potassium	627	U	---	702	U	---	745	U	---	709	U	---	651	U	---	741	U	---	733	U	---
Selenium	0.21	UN	UJ	0.24	BN	J	0.25	UN	UJ	0.24	UN	UJ	0.22	UN	UJ	0.25	UN	UJ	0.25	UN	UJ
Silver	0.74	U	---	0.82	U	---	0.87	U	---	0.83	U	---	0.76	U	---	0.87	U	---	0.86	U	---
Sodium	51.7	B	U	53.7	B	U	82.8	B	U	48.0	B	U	89.6	B	U	70.4	B	U	135.0	B	U
Thallium	0.21	UW	UJ	0.24	U	---	0.25	U	---	0.24	U	---	0.22	U	---	0.25	U	---	0.25	U	---
Vanadium	13.8		---	22.2		---	34.0		---	27.4		---	13.2		---	14.9		---	21.8		---
Zinc	12.5		---	9.5		---	17.9		---	10.4		---	10.4		---	13.6		---	14.7		---
Cyanide	0.12	U	---	0.14	U	---	0.16	B	J	0.14	U	---	0.15	B	J	0.59	B	J	0.15	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**	SA3-10D-SS-D	Q*	DV**	SA3-10E-SS	Q*	DV**	SA4-11A-SS	Q*	DV**
Aluminum	5570		---	3900		---	2280		---	3100		---	3580		NP	4370		---	2190		---
Antimony	13.3	U	---	14.1	U	---	13.6	U	---	13.3	U	---	13.3	U	NP	14.4	U	---	12.4	UN	J
Arsenic	1.30	BN	J	1.80	BN	J	1.00	BN	J	1.20	BN	J	1.20	B	NP	1.70	BN	J	1.90	B	J
Barium	18.9	B	J	22.1	B	J	19.0	B	J	42.7	B	J	43.6	B	NP	35.3	B	J	17.8	B	J
Beryllium	0.26	B	U	0.14	B	U	0.13	B	U	0.25	B	U	0.13	B	NP	0.27	B	U	0.06	U	---
Cadmium	0.93	B	U	0.75	U	---	0.96	B	U	0.70	U	---	0.70	U	NP	1.30		---	0.65	U	---
Calcium	723	B	J	235	B	J	351	B	J	436	B	J	437	B	NP	446	B	J	612	B	J
Chromium	7.6		J	5.0		J	3.0		J	7.7		J	4.6		NP	4.2		J	5.3		---
Cobalt	3.30	B	J	1.60	B	J	1.00	B	J	2.20	B	J	2.10	B	NP	3.10	B	J	2.90	B	U
Copper	2.50	B	J	0.63	U	UJ	0.60	U	---	2.30	B	J	2.30	B	NP	1.90	B	J	3.80	B	U
Iron	10900		---	4360		---	2790		---	3450		---	3370		NP	4130		---	3260		---
Lead	6.6	N	J	9.3	N	J	7.5	N	J	10.0	N	J	9.7		NP	9.6	N	J	11.5		J
Magnesium	376.0	B	J	196.0	B	J	124.0	B	J	157.0	B	J	173.0	B	NP	212.0	B	J	101.0	B	J
Manganese	107.0		---	56.5		---	104.0		---	350.0		---	332.0		NP	233.0		---	41.3		---
Mercury	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	NP	0.06	U	---	0.14		---
Nickel	4.7	U	---	5.0	U	---	4.8	U	---	4.7	U	---	4.7	U	NP	5.1	U	---	4.4	U	---
Potassium	676	U	---	719	U	---	774	B	J	676	U	---	676	U	NP	732	U	---	630	U	---
Selenium	0.23	UN	R	0.24	UN	R	0.23	UN	R	0.23	UWN	R	0.23	U	NP	0.25	UN	R	0.21	U	---
Silver	0.79	U	UJ	0.84	U	---	0.81	U	---	0.79	U	---	0.79	U	NP	0.86	U	---	0.74	U	---
Sodium	60.4	B	U	56.2	B	U	43.9	B	U	46.9	B	U	60.5	B	NP	59.7	B	U	42.5	B	U
Thallium	0.23	UN	UJ	0.24	UN	UJ	0.23	UN	UJ	0.23	UN	UJ	0.23	U	NP	0.25	UN	UJ	0.21	U	---
Vanadium	18.7		---	9.7	B	J	5.9	B	J	7.7	B	J	7.5	B	NP	9.5	B	J	5.4	B	J
Zinc	4.8		---	4.6	B	J	2.4	B	J	4.3	B	J	4.3	B	NP	4.5	B	J	13.3		---
Cyanide	0.14	U	---	0.15	U	---	0.13	U	---	0.14	U	---	0.15	U	NP	0.14	U	---	0.14	U	---

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**	SA4-12A-SS-D	Q*	DV**	SA4-12B-SS	Q*	DV**
Aluminum	3380		---	3810		---	4350		---	3070		---	14200		---	17000		NP	7840		---
Antimony	12.8	UN	J	12.4	UN	J	13.3	UN	J	12.5	UN	J	14.8	N	R	14.1	U	NP	13.2	BN	R
Arsenic	1.80	B	J	1.90	BW	J	4.00		---	1.70	B	J	1.60	B	J	2.20	B	NP	2.20		---
Barium	32.9	B	J	48.2		---	39.2	B	J	15.8	B	J	22.7	B	J	24.6	B	NP	49.3		---
Beryllium	0.07	U	---	0.13	B	J	0.07	U	---	0.06	U	---	0.23	B	J	0.15	B	NP	0.07	U	---
Cadmium	0.67	U	---	0.66	U	---	0.70	U	---	0.66	U	---	0.90	B	J	0.75	U	NP	0.69	U	---
Calcium	712	B	J	2150		---	1490		---	709	B	J	1170	B	J	1170	B	NP	576	B	J
Chromium	6.6		---	15.9		---	5.3		---	6.6		---	14.8		---	17.0		NP	7.9		---
Cobalt	2.50	B	U	3.60	B	U	3.80	B	U	1.80	B	U	4.90	B	J	3.40	B	NP	3.80	B	J
Copper	4.80	B	J	11.10		---	17.70		---	4.90	B	J	5.40	B	J	3.70	B	NP	4.50	B	J
Iron	4220		---	10600		---	7280		---	4540		---	19700		---	17500		NP	12300		---
Lead	8.8		J	54.3		J	22.9		J	6.6		J	6.9		---	7.8		NP	9.1		---
Magnesium	171.0	B	J	190.0	B	J	251.0	B	J	161.0	B	U	1300.0		---	1500.0		NP	773.0	B	J
Manganese	40.9		---	94.9		---	104.0		---	26.4		---	24.8		---	28.9		NP	16.1		---
Mercury	0.21		---	0.05	U	---	0.07	B	U	0.06	B	U	0.06	U	---	0.06	U	NP	0.06	U	---
Nickel	4.8	B	J	4.4	U	---	4.7	U	---	4.4	U	---	5.0	U	---	5.0	U	NP	4.7	U	---
Potassium	648	U	---	633	U	---	676	U	---	633	U	---	717	U	---	717	U	NP	667	U	---
Selenium	0.22	U	---	0.21	U	---	0.23	U	---	0.21	U	---	0.69	BN	J	0.34	B	NP	0.22	UN	UJ
Silver	0.76	U	---	0.74	U	---	0.79	U	---	0.74	U	---	0.84	U	---	0.84	U	NP	0.78	U	---
Sodium	47.5	B	U	48.2	B	U	69.9	B	U	50.1	B	U	56.6	B	J	67.4	B	NP	44.7	B	J
Thallium	0.22	U	---	0.21	U	---	0.23	U	---	0.21	U	---	0.24	U	---	0.24	U	NP	0.22	U	---
Vanadium	7.0	B	J	9.7	B	J	10.3	B	J	9.5	B	J	29.1		---	26.5		NP	17.2		---
Zinc	33.1		---	29.3		---	50.9		---	6.7		---	13.5		---	16.8		NP	10.2		---
Cyanide	0.14	U	---	0.14	U	---	0.15	U	---	0.49	B	J	0.19	B	J	0.16	U	NP	0.15	U	---

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	Table A-5.1 Metals/Cyanide																							
	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**	SA4-13D-SS	Q*	DV**
Aluminum	3750		---	4460		---	6490		---	2710		---	3880		J	6830		J	6520		J
Antimony	13.5	UN	R	13.6	UN	R	12.4	UN	R	12.2	UN	UJ	12.6	UN	UJ	12.3	UN	UJ	13.0	UN	UJ
Arsenic	0.73	B	J	0.64	B	J	1.40	B	J	1.10	B	J	1.50	B	J	1.80	B	J	1.50	B	J
Barium	429.0		---	32.6	B	J	19.5	B	J	30.3	B	J	25.3	B	J	45.0		---	35.2	B	J
Beryllium	0.07	U	---	0.07	U	---	0.06	U	---	0.06	B	U	0.06	U	---	0.13	B	U	0.13	B	U
Cadmium	0.71	U	---	0.78	B	J	0.65	U	---	0.64	U	---	0.66	U	---	0.65	U	---	0.69	U	---
Calcium	406	B	J	361	B	J	403	B	J	543	B	J	657	B	J	5020		---	393	B	J
Chromium	4.3		---	5.4		---	12.2		---	9.6		---	6.8		---	28.9		---	5.2		---
Cobalt	0.96	U	---	2.10	B	J	2.20	B	J	2.90	B	J	2.50	B	J	3.90	B	J	2.30	B	J
Copper	0.60	U	---	1.20	B	J	1.90	B	J	4.80	B	U	3.80	B	U	6.50		---	4.30	B	U
Iron	3530		---	3700		---	7420		---	4000		J	4400		J	8210		J	5420		J
Lead	9.8		---	7.9		---	11.1		---	12.9	N	J	13.4	SN	J	16.6	N	J	10.1	N	J
Magnesium	197.0	B	J	223.0	B	J	271.0	B	J	214.0	B	J	206.0	B	J	445.0	B	J	331.0	B	J
Manganese	143.0		---	116.0		---	32.6		---	44.2		J	37.0		J	82.5		J	98.8		J
Mercury	0.06	U	---	0.06	U	---	0.05	U	---	0.05	U	---	0.05	U	---	0.06	B	J	0.06	U	---
Nickel	4.8	U	---	4.9	U	---	4.4	U	---	4.3	U	---	4.5	U	---	19.1		---	4.6	U	---
Potassium	685	U	---	693	U	---	629	U	---	618	U	---	639	U	---	624	U	---	661	U	---
Selenium	0.29	BN	J	0.23	UWN	UJ	0.21	UN	UJ	0.21	U	---	0.21	U	---	0.21	U	---	0.22	U	---
Silver	0.80	U	---	0.81	U	---	0.74	U	---	0.73	U	---	0.75	U	---	0.73	U	---	0.78	U	---
Sodium	31.9	B	J	25.2	B	U	28.3	B	U	41.7	B	U	49.3	B	U	80.1	B	J	43.6	B	U
Thallium	0.23	U	---	0.23	U	---	0.21	U	---	0.21	U	---	0.21	U	---	0.21	U	---	0.22	U	---
Vanadium	7.5	B	J	8.5	B	J	14.2		---	7.4	B	J	9.9	B	J	14.4		---	11.8		---
Zinc	6.6		---	7.2		---	9.8		---	9.8		---	12.9		---	19.3		---	90.9		---
Cyanide	0.68		---	0.15	U	---	0.14	U	---	0.13	U	---	0.14	U	---	0.14	U	---	0.14	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14A-SS-D	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**
Aluminum	2720		J	3730		---	3680		NP	7090		---	2980		---	2720		---	5240		---
Antimony	13.7	UN	UJ	12.4	UN	R	12.4	UN	NP	13.0	UN	R	12.8	UN	R	12.2	UN	R	12.5	UN	R
Arsenic	0.69	B	U	0.67	B	J	1.01	B	NP	1.60	B	J	1.10	B	J	1.00	B	J	1.00	B	J
Barium	38.6	B	J	18.4	B	J	15.2	B	NP	17.2	B	J	11.5	B	J	10.0	B	J	15.3	B	J
Beryllium	0.14	B	U	0.06	U	---	0.06	U	NP	0.07	U	---	0.07	U	---	0.06	U	---	0.06	U	---
Cadmium	0.73	U	---	0.65	U	---	1.01	B	NP	0.69	U	---	0.72	B	J	0.65	U	---	0.66	U	---
Calcium	755	B	J	329	B	J	342	B	NP	215	B	J	283	B	J	202	B	J	199	B	J
Chromium	3.0		---	4.2		---	3.8		NP	7.6		---	4.5		---	3.9		---	5.5		---
Cobalt	1.70	B	J	0.89	U	---	0.90	U	NP	2.00	B	J	4.40	B	J	2.20	B	J	1.80	B	J
Copper	1.40	B	U	1.10	B	U	1.10	B	NP	2.00	B	U	2.70	B	U	4.90	B	U	6.30	[illegible]	---
Iron	2370		J	5560		---	4940		NP	6660		---	4350		---	4480		---	5530		---
Lead	7.6	SN	J	4.2		---	5.3		NP	7.9		---	10.2		---	78.9	S	---	6.1		---
Magnesium	186.0	B	J	173.0	B	J	173.0	B	NP	237.0	B	J	104.0	B	J	96.6	B	J	162.0	B	J
Manganese	83.3		J	54.1		---	70.5		NP	19.5		---	29.5		---	50.6		---	21.1		---
Mercury	0.06	U	---	0.05	U	---	0.05	U	NP	0.06	U	---	0.05	U	---	0.05	U	---	0.05	U	---
Nickel	4.9	U	---	4.4	U	---	4.4	U	NP	4.6	U	---	4.6	U	---	4.4	U	---	4.4	U	---
Potassium	699	U	---	629	U	---	629	U	NP	662	U	---	650	U	---	622	U	---	633	U	---
Selenium	0.23	U	---	0.24	BN	UJ	0.21	U	NP	0.24	BN	UJ	0.22	UN	UJ	0.21	UN	J	0.21	UN	UJ
Silver	0.82	U	---	0.74	U	---	0.74	U	NP	0.78	U	---	0.76	U	---	0.73	U	---	0.74	U	---
Sodium	47.2	B	U	15.4	B	U	16.4	B	NP	24.6	B	U	25.6	B	U	11.3	B	U	18.5	B	U
Thallium	0.23	U	---	0.21	U	---	0.21	U	NP	0.22	U	---	0.22	U	---	0.21	U	---	0.21	U	---
Vanadium	4.8	B	J	11.2		---	11.1		NP	18.1		---	10.3	B	J	10.2	B	J	13.3		---
Zinc	7.0		---	4.4		---	5.0		NP	6.3		---	5.7		---	5.2		---	4.7		---
Cyanide	0.15	U	---	0.14	U	---	0.14	U	NP	0.15	U	---	0.14	U	---	0.14	U	---	0.14	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA4-15A-SS	Q*	DV**	SA4-15A-SS-D	Q*	DV**	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**
Aluminum	4630		---	4430		NP	4420		---	7340		---	5250		---	7240		---	5620		---
Antimony	12.2	U	---	12.2	U	NP	14.0	U	---	23.6		---	13.6	U	---	13.7	U	---	13.0	UN	UJ
Arsenic	1.50	B	J	2.00	B	NP	3.80		---	2.70		---	5.00		---	3.40		---	0.99	BW	J
Barium	19.1	B	J	18.1	B	NP	33.3	B	J	29.6	B	J	75.8		---	28.6	B	J	48.3		---
Beryllium	0.13	B	U	0.07	B	NP	0.23	B	U	0.48	B	U	0.36	B	U	0.51	B	U	0.07	U	---
Cadmium	0.65	U	---	0.65	U	NP	0.74	U	---	0.69	U	---	0.72	U	---	0.72	U	---	0.69	U	---
Calcium	501	B	J	520	B	NP	454	B	J	937	B	J	2790		---	658	B	J	318	BE	J
Chromium	4.6		---	4.6		NP	6.4		---	9.0		---	8.3		---	10.1		---	4.3		---
Cobalt	2.00	B	U	2.50	B	NP	2.70	B	U	4.30	B	U	4.50	B	U	3.00	B	U	3.70	B	J
Copper	39.00	E	J	25.23		NP	50.70	E	J	9950.00	E	J	73.90	E	J	84.70	E	J	0.58	UN	---
Iron	6440		---	4880		NP	5910		---	9220		---	17700		---	12000		---	5020		---
Lead	5.5	S	---	5.9		NP	34.6		---	120.0		---	16.8	S	---	10.1	S	---	6.8		---
Magnesium	245.0	B	J	249.0	B	NP	273.0	B	J	621.0	B	J	300.0	B	J	391.0	B	J	259.0	B	J
Manganese	40.1		J	20.0		NP	198.0		J	21.4		J	127.0		J	48.9		J	468.0		---
Mercury	0.05	U	---	0.05	U	NP	0.07	B	J	0.06	U	---	0.06	U	---	0.06	B	J	0.08	B	J
Nickel	4.4	U	---	4.4	U	NP	5.0	U	---	4.7	B	J	7.3	B	J	4.9	U	-	4.6	U	---
Potassium	621	U	---	621	U	NP	709	U	---	659	U	---	690	U	---	695	U	---	660	U	---
Selenium	0.21	UN	UJ	0.21	U	NP	0.24	UN	UJ	0.22	UN	UJ	0.23	UN	UJ	0.23	UN	UJ	0.22	U	---
Silver	0.73	U	---	0.73	U	NP	0.83	U	---	0.77	U	---	0.81	U	---	0.82	U	---	0.77	U	---
Sodium	52.2	B	U	43.9	B	NP	45.4	B	U	67.5	B	U	57.9	B	U	104.0	B	J	50.5	B	U
Thallium	0.21	U	---	0.21	U	NP	0.24	U	---	0.22	U	---	0.23	U	---	0.23	U	---	0.22	U	---
Vanadium	12.5		---	10.3	B	NP	13.1		---	17.7		---	15.4		---	24.6		---	9.9	B	J
Zinc	8.0		---	6.1		NP	11.4		---	131.0		---	114.0		---	18.8		---	8.3	N	UJ
Cyanide	0.14	U	---	0.14	U	NP	0.16	U	---	0.19	B	J	0.15	U	---	0.22	B	J	0.15	U	---

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA5-16B-SS	Q*	DV**	SA5-16C-SS	Q*	DV**	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**
Aluminum	2310		---	7520		---	2250		---	3850		---	3920		---	5540		---	6370		---
Antimony	12.8	UN	UJ	12.8	UN	UJ	12.5	UN	UJ	15.2	UN	UJ	13.3	UN	R	13.3	UN	R	12.8	UN	R
Arsenic	0.74	BW	J	1.50	BW	J	5.70	S	---	2.60	W	J	0.86	B	J	1.10	B	J	1.60	B	J
Barium	28.9	B	J	22.6	B	J	7.2	B	J	38.5	B	J	68.8		---	60.0		---	69.4		---
Beryllium	0.07	U	---	0.07	U	---	0.06	U	---	0.08	U	---	0.07	U	---	0.07	U	---	0.07	B	J
Cadmium	0.68	U	---	0.68	U	---	0.66	U	---	0.80	U	---	0.70	U	---	0.70	U	---	0.86	B	J
Calcium	307	BE	J	620	BE	J	295	BE	J	9480	E	J	746	B	J	738	B	J	5500		---
Chromium	3.1		---	24.3		---	14.8		---	4.3		---	2.5		---	4.3		---	5.5		---
Cobalt	1.90	B	J	2.80	B	U	2.30	B	U	3.50	B	U	3.20	B	J	4.90	B	J	4.10	B	J
Copper	0.57	UN	---	4.30	BN	J	0.55	UN	---	0.67	UN	---	0.92	B	U	0.95	B	U	5.30	B	J
Iron	3290		---	10400		---	7140		---	5570		---	3150		---	4540		---	7020		---
Lead	5.8	S	---	6.9		---	12.4		---	11.7		---	10.5		---	11.9		---	36.3		---
Magnesium	138.0	B	J	251.0	B	J	101.0	B	J	553.0	B	J	178.0	B	J	259.0	B	J	389.0	B	J
Manganese	267.0		---	38.3		---	44.7		---	147.0		---	620.0		---	669.0		---	434.0		---
Mercury	0.07	B	J	0.05	U	---	0.05	U	---	0.06	U	---	0.08	B	J	0.06	U	---	0.05	U	---
Nickel	4.6	U	---	4.6	U	---	4.4	U	---	5.4	U	---	4.7	U	---	4.8	U	---	4.6	U	---
Potassium	651	U	---	651	U	---	633	U	---	772	U	---	676	U	---	678	U	---	651	U	---
Selenium	0.22	UW	UJ	0.22	UW	UJ	0.21	UW	UJ	0.65	BW	J	0.23	UWN	UJ	0.31	BN	UJ	0.25	BN	UJ
Silver	0.76	U	---	0.76	U	---	0.74	U	---	0.91	U	---	0.79	U	---	0.80	U	---	0.76	U	---
Sodium	40.7	B	U	53.5	B	U	54.7	B	U	64.4	B	U	29.3	B	U	25.2	B	U	26.2	B	U
Thallium	0.22	U	---	0.28	B	J	0.21	U	---	0.26	U	---	0.23	U	---	0.23	U	---	0.22	U	---
Vanadium	5.6	B	J	20.8		---	17.0		---	8.6	B	J	7.0	B	J	9.4	B	J	13.2		---
Zinc	8.4	N	UJ	9.5	N	UJ	4.8	N	UJ	17.5	N	UJ	5.2		---	7.5		---	24.4		---
Cyanide	0.14	U	---	0.15	B	J	0.14	U	---	0.17	U	---	0.15	U	---	0.15	U	---	0.17	B	J

* Data Qualifier. ** Data Validation.

Table A-5.1 Metals/Cyanide

Compound (mg/Kg)	SA5-17D-SS	Q*	DV**	SA5-17E-SS	Q*	DV**	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**
Aluminum	6930		---	6660		---	3130		---	4600		---	2630		---	2750		---	4450		---
Antimony	13.7	UN	R	13.4	UN	R	13.3	UN	UJ	13.2	UN	UJ	13.1	UN	UJ	12.9	UN	UJ	13.1	UN	UJ
Arsenic	1.20	B	J	0.93	B	J	1.10	BW	---	1.40	BW	J	0.94	BW	J	0.91	BW	J	1.30	BW	J
Barium	76.7		---	107.0		---	22.4	B	J	19.9	B	J	13.5	B	J	19.7	B	J	26.2	B	J
Beryllium	0.07	U	---	0.27	B	J	0.07	U	---	0.07	U	---	0.07	U	---	0.07	U	---	0.07	U	---
Cadmium	0.72	U	---	0.71	U	---	0.70	U	---	0.70	U	---	0.69	U	---	0.68	U	---	0.69	U	---
Calcium	842	B	J	775	B	J	263	BE	J	301	BE	J	235	BE	J	287	BE	J	287	BE	J
Chromium	4.9		---	5.2		---	6.9		---	5.2		---	3.8		---	2.6		---	4.0		---
Cobalt	4.00	B	J	5.20	B	J	2.50	B	U	1.70	B	U	0.94	U	---	2.50	B	U	2.10	B	U
Copper	0.61	U	---	1.20	B	U	0.59	UN	---	0.59	UN	---	0.58	UN	---	0.57	UN	---	0.58	UN	---
Iron	4740		---	4880		---	3430		---	4520		---	3210		---	2500		---	3880		---
Lead	9.2		---	63.6		---	19.7		---	16.1		---	23.1		---	6.4		---	6.7		---
Magnesium	343.0	B	J	307.0	B	J	151.0	B	J	209.0	B	J	133.0	B	J	140.0	B	J	170.0	B	J
Manganese	754.0		---	1150.0		---	207.0		---	72.5		---	43.8		---	144.0		---	157.0		---
Mercury	0.06	B	J	0.06	U	---	0.06	B	J	0.06	U	---	0.06	U	---	0.05	U	-	0.07	B	J
Nickel	4.9	U	---	4.8	U	---	4.7	U	---	4.7	U	---	4.7	U	---	4.6	U	---	4.6	U	---
Potassium	696	U	---	681	U	---	674	U	---	726	B	J	665	U	---	653	U	---	725	B	J
Selenium	0.37	BN	UJ	0.44	BN	UJ	0.23	UW	UJ	0.23	UW	UJ	0.22	U	---	0.22	UW	UJ	0.22	UW	UJ
Silver	0.82	U	---	0.80	U	---	0.79	U	---	0.79	U	---	0.78	U	---	0.77	U	---	0.78	U	---
Sodium	28.0	B	U	25.3	B	U	57.2	B	U	50.0	B	U	41.5	B	U	52.7	B	U	50.7	B	U
Thallium	0.23	U	---	0.23	U	---	0.23	UW	UJ	0.23	UW	UJ	0.22	UW	UJ	0.22	U	---	0.22	U	---
Vanadium	10.9	B	J	10.5	B	J	6.7	B	J	8.6	B	J	5.3	B	J	5.5	B	J	8.9	B	J
Zinc	8.5		---	9.1		---	7.6	N	UJ	8.2	N	UJ	10.6	N	UJ	4.8	N	UJ	6.9	N	UJ
Cyanide	0.15	U	---	0.15	U	---	0.15	U	---	0.15	U	---	0.15	U	---	0.14	U	---	0.15	U	---

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	Table A-5.1 Metals/Cyanide														
	REFC-19A-SS	Q*	DV**	REFC-19B-SS	Q*	DV**	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**
Aluminum	3490		---	5010		---	14600		---	1980		---	3820		---
Antimony	12.6	UN	UJ	13.5	UN	UJ	16.2	UN	UJ	12.1	UN	UJ	12.5	UN	UJ
Arsenic	1.20	BW	J	2.00	BW	J	2.70	B	J	0.66	BW	J	1.10	BW	J
Barium	21.5	B	J	36.4	B	J	80.6		---	9.6	B	J	40.0	B	J
Beryllium	0.06	U	---	0.07	U	---	0.33	B	J	0.06	U	---	0.06	U	---
Cadmium	0.67	U	---	0.71	U	---	0.85	U	---	0.64	U	---	0.66	U	---
Calcium	306	BE	J	519	BE	J	671	BE	J	273	BE	J	539	BE	J
Chromium	2.4		---	4.4		---	10.4		---	0.95	U	---	2.2		---
Cobalt	0.90	U	---	2.00	B	U	5.60	B	J	1.10	B	J	2.10	B	J
Copper	10.70	N	J	0.60	UN	---	0.72	UN	---	0.54	UN	---	0.55	UN	---
Iron	3540		---	6700		---	11700		---	1780		---	3340		---
Lead	4.6		---	7.7		---	12.6		---	1.8		---	5.6		---
Magnesium	186.0	B	J	210.0	B	J	582.0	B	J	120.0	B	J	218.0	B	J
Manganese	74.8		---	211.0		---	385.0		---	60.5		---	241.0		---
Mercury	0.05	U	---	0.06	U	---	0.07	U	---	0.05	U	---	0.05	U	---
Nickel	4.5	U	---	4.8	U	---	7.8	B	J	4.3	U	---	4.5	U	---
Potassium	641	U	---	687	U	---	822	U	---	617	U	---	656	B	J
Selenium	0.22	UW	UJ	0.23	UW	UJ	0.28	UW	UJ	0.21	U	---	0.21	U	---
Silver	0.75	U	---	0.81	U	---	0.96	U	---	0.72	U	---	0.75	U	---
Sodium	49.1	B	U	52.6	B	U	62.8	B	U	87.9	B	U	117.0	B	J
Thallium	0.22	U	---	0.23	U	---	0.28	U	---	0.21	U	---	0.21	U	---
Vanadium	7.7	B	J	14.6		---	25.6		---	4.1	B	J	6.9	B	J
Zinc	7.2	N	UJ	16.7	N	UJ	24.4	N	J	4.7	N	J	7.4	N	J
Cyanide	0.14	U	---	0.15	U	---	0.27	B	J	0.16	B	J	0.17	B	J

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA1-1A-SS	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1E-SS	Q*	DV**	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**
Alpha-BHC	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Beta-BHC	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Delta-BHC	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Gamma-BHC	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Heptachlor	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Aldrin	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Heptachlor epoxide	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Endosulfan I	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---	18	U	---	18	U	---
Dieldrin	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
4,4'-DDE	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
Endrin	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
Endosulfan II	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
4,4'-DDD	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
Endosulfan sulfate	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
4,4'-DDT	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
Methoxychlor	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---
Endrin ketone	36	U	---	34	U	---	34	U	---	35	U	---	34	U	---	35	U	---	37	U	---
Alpha-chlordane	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---
Gamma-chlordane	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---
Toxaphene	360	U	---	340	U	---	340	U	---	350	U	---	340	U	---	350	U	---	370	U	---
Aroclor-1016	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---
Aroclor-1221	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---
Aroclor-1232	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---
Aroclor-1242	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA1-1A-SS	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1E-SS	Q*	DV**	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**			
Aroclor-1248	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---	180	U	---	180	U	---			
Aroclor-1254	360	U	---	340	U	---	340	U	---	350	U	---	340	U	---	350	U	---	370	U	---			
Aroclor-1260	360	U	---	340	U	---	340	U	---	350	U	---	340	U	---	350	U	---	370	U	---			

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**
Alpha-BHC	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Beta-BHC	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Delta-BHC	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Gamma-BHC	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Heptachlor	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Aldrin	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Heptachlor epoxide	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Endosulfan I	17	U	---	19	U	---	18	U	---	20	U	---	17	U	---	17	U	---	17	U	---
Dieldrin	34	U	---	38	U	---	37	U	---	41	U	---	34	U	---	34	U	---	35	U	---
4,4'-DDE	34	U	---	38	U	---	37	U	---	210		---	34	U	---	34	U	---	35	U	---
Endrin	34	U	---	38	U	---	37	U	---	41	U	---	34	U	---	34	U	---	35	U	---
Endosulfan II	34	U	---	38	U	---	37	U	---	41	U	---	34	U	---	34	U	---	35	U	---
4,4'-DDD	34	U	---	38	U	---	37	U	---	41		---	34	U	---	34	U	---	35	U	---
Endosulfan sulfate	34	U	---	38	U	---	37	U	---	41	U	---	34	U	---	34	U	---	35	U	---
4,4'-DDT	34	U	---	38	U	---	37	U	---	190		---	34	U	---	34	U	---	35	U	---
Methoxychlor	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---
Endrin ketone	34	U	---	38	U	---	37	U	---	41	U	---	34	U	---	34	U	---	35	U	---
Alpha-chlordane	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---
Gamma-chlordane	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---
Toxaphene	340	U	---	380	U	---	370	U	---	410	U	---	340	U	---	340	U	---	350	U	---
Aroclor-1016	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1221	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1232	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1242	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**
	Table A-5.2 Pesticides/PCB's																				
Aroclor-1248	170	U	---	190	U	---	180	U	---	200	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1254	340	U	---	380	U	---	370	U	---	410	U	---	340	U	---	340	U	---	350	U	---
Aroclor-1260	340	U	---	380	U	---	370	U	---	410	U	---	340	U	---	340	U	---	350	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA1-4A-SS	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**	SA2-5B-SS	Q*	DV**
Alpha-BHC	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Beta-BHC	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Delta-BHC	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Gamma-BHC	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Heptachlor	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Aldrin	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Heptachlor epoxide	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Endosulfan I	20	U	---	19	U	---	17	U	---	19	U	---	20	U	---	19	U	---	18	U	---
Dieldrin	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
4,4'-DDE	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
Endrin	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
Endosulfan II	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
4,4'-DDD	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
Endosulfan sulfate	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
4,4'-DDT	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
Methoxychlor	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---
Endrin ketone	40	U	---	37	U	---	34	U	---	38	U	---	40	U	---	37	U	---	37	U	---
Alpha-chlordane	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---
Gamma-chlordane	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---
Toxaphene	400	U	---	370	U	---	340	U	---	380	U	---	400	U	---	370	U	---	370	U	---
Aroclor-1016	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---
Aroclor-1221	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---
Aroclor-1232	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---
Aroclor-1242	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA1-4A-SS	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**	SA2-5B-SS	Q*	DV**
Aroclor-1248	200	U	---	190	U	---	170	U	---	190	U	---	200	U	---	190	U	---	180	U	---
Aroclor-1254	400	U	---	370	U	---	340	U	---	380	U	---	400	U	---	370	U	---	370	U	---
Aroclor-1260	400	U	---	370	U	---	340	U	---	380	U	---	400	U	---	370	U	---	370	U	---

Compound (ug/Kg)	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**	SA2-6D-SS	Q*	DV**
	Table A-5.2 Pesticides/PCB's																				
Alpha-BHC	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Beta-BHC	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Delta-BHC	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Gamma-BHC	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Heptachlor	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Aldrin	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Heptachlor epoxide	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Endosulfan I	19	U	---	17	U	---	19	U	---	19	U	---	17	U	---	17	U	---	17	U	---
Dieldrin	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
4,4'-DDE	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
Endrin	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
Endosulfan II	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
4,4'-DDD	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
Endosulfan sulfate	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
4,4'-DDT	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
Methoxychlor	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---
Endrin ketone	38	U	---	35	U	---	37	U	---	37	U	---	34	U	---	33	U	---	35	U	---
Alpha-chlordane	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---
Gamma-chlordane	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---
Toxaphene	380	U	---	350	U	---	370	U	---	370	U	---	340	U	---	330	U	---	350	U	---
Aroclor-1016	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1221	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1232	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1242	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**	SA2-6D-SS	Q*	DV**
Aroclor-1248	190	U	---	170	U	---	190	U	---	190	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1254	380	U	---	350	U	---	370	U	---	370	U	---	340	U	---	330	U	---	350	U	---
Aroclor-1260	380	U	---	350	U	---	370	U	---	370	U	---	340	U	---	330	U	---	350	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**	SA3-8A-SS	Q*	DV**
Alpha-BHC	18	U	---	19	U	---	19	U	---	17	U	---	18	U	---	17	U	---	19	U	---
Beta-BHC	18	U	---	19	U	---	19	U	---	17	U	---	18	U	---	17	U	---	19	U	---
Delta-BHC	18	U	---	19	U	---	19	U	---	17	U	---	18	U	---	17	U	---	19	U	---
Gamma-BHC	18	U	---	19	U	---	19	U	---	17	U	---	18	U	---	17	U	---	19	U	---
Heptachlor	18	U	---	19	U	---	19	U	---	17	U	---	18	U	---	17	U	---	19	U	---
Aldrin	18	U	---	19	U	---	19	U	---	17	U	---	18	U	---	17	U	---	19	U	---
Heptachlor epoxide	18	U	---	19	U	---	19	U	---	17	U	---	62		---	17	U	---	19	U	---
Endosulfan I	18	U	---	19	U	---	19	U	---	17	U	---	18	U	---	17	U	---	19	U	---
Dieldrin	35	U	---	37	U	---	39	U	---	42		---	320		---	34	U	---	38	U	---
4,4'-DDE	35	U	---	37	U	---	39	U	---	35	U	---	35	U	---	34	U	---	38	U	---
Endrin	35	U	---	37	U	---	39	U	---	35	U	---	35	U	---	34	U	---	38	U	---
Endosulfan II	35	U	---	37	U	---	39	U	---	35	U	---	38		---	34	U	---	38	U	---
4,4'-DDD	35	U	---	37	U	---	39	U	---	35	U	---	35	U	---	34	U	---	38	U	---
Endosulfan sulfate	35	U	---	37	U	---	39	U	---	35	U	---	35	U	---	34	U	---	38	U	---
4,4'-DDT	35	U	---	37	U	---	39	U	---	35	U	---	35	U	---	34	U	---	38	U	---
Methoxychlor	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---
Endrin ketone	35	U	---	37	U	---	39	U	---	35	U	---	35	U	---	34	U	---	38	U	---
Alpha-chlordane	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---
Gamma-chlordane	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---
Toxaphene	350	U	---	370	U	---	390	U	---	350	U	---	350	U	---	340	U	---	380	U	---
Aroclor-1016	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---
Aroclor-1221	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---
Aroclor-1232	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---
Aroclor-1242	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**	SA3-8A-SS	Q*	DV**
Aroclor-1248	180	U	---	190	U	---	190	U	---	170	U	---	180	U	---	170	U	---	190	U	---
Aroclor-1254	350	U	---	370	U	---	390	U	---	350	U	---	350	U	---	340	U	---	380	U	---
Aroclor-1260	350	U	---	370	U	---	390	U	---	350	U	---	350	U	---	340	U	---	380	U	---

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**	SA3-9C-SS	Q*	DV**
Alpha-BHC	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Beta-BHC	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Delta-BHC	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Gamma-BHC	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Heptachlor	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Aldrin	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Heptachlor epoxide	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Endosulfan I	17	U	---	16	U	---	17	U	---	19	U	---	20	U	---	19	U	---	17	U	---
Dieldrin	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
4,4'-DDE	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
Endrin	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
Endosulfan II	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
4,4'-DDD	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
Endosulfan sulfate	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
4,4'-DDT	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
Methoxychlor	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---
Endrin ketone	34	U	---	33	U	---	34	U	---	38	U	---	40	U	---	38	U	---	35	U	---
Alpha-chlordane	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---
Gamma-chlordane	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---
Toxaphene	340	U	---	330	U	---	340	U	---	380	U	---	400	U	---	380	U	---	350	U	---
Aroclor-1016	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---
Aroclor-1221	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---
Aroclor-1232	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---
Aroclor-1242	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**	SA3-9C-SS	Q*	DV**
Aroclor-1248	170	U	---	160	U	---	170	U	---	190	U	---	200	U	---	190	U	---	170	U	---
Aroclor-1254	340	U	---	330	U	---	340	U	---	380	U	---	400	U	---	380	U	---	350	U	---
Aroclor-1260	340	U	---	330	U	---	340	U	---	380	U	---	400	U	---	380	U	---	350	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**	SA3-10E-SS	Q*	DV**
Alpha-BHC	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Beta-BHC	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Delta-BHC	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Gamma-BHC	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Heptachlor	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Aldrin	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Heptachlor epoxide	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Endosulfan I	20	U	---	20	U	---	18	U	---	19	U	---	18	U	---	18	U	---	19	U	---
Dieldrin	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
4,4'-DDE	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
Endrin	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
Endosulfan II	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
4,4'-DDD	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
Endosulfan sulfate	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
4,4'-DDT	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
Methoxychlor	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---
Endrin ketone	40	U	---	39	U	---	36	U	---	39	U	---	37	U	---	36	U	---	39	U	---
Alpha-chlordane	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---
Gamma-chlordane	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---
Toxaphene	400	U	---	390	U	---	360	U	---	390	U	---	370	U	---	360	U	---	390	U	---
Aroclor-1016	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---
Aroclor-1221	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---
Aroclor-1232	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---
Aroclor-1242	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**	SA3-10E-SS	Q*	DV**
Aroclor-1248	200	U	---	200	U	---	180	U	---	190	U	---	180	U	---	180	U	---	190	U	---
Aroclor-1254	400	U	---	390	U	---	360	U	---	390	U	---	370	U	---	360	U	---	390	U	---
Aroclor-1260	400	U	---	390	U	---	360	U	---	390	U	---	370	U	---	360	U	---	390	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA4-11A-SS	Q*	DV**	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**	SA4-12B-SS	Q*	DV**
Alpha-BHC	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Beta-BHC	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Delta-BHC	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Gamma-BHC	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Heptachlor	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Aldrin	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Heptachlor epoxide	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Endosulfan I	16	U	---	16	U	---	16	U	---	17	U	---	16	U	---	19	U	---	18	U	---
Dieldrin	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
4,4'-DDE	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
Endrin	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
Endosulfan II	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
4,4'-DDD	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
Endosulfan sulfate	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
4,4'-DDT	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
Methoxychlor	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---
Endrin ketone	33	U	---	32	U	---	32	U	---	34	U	---	32	U	---	38	U	---	35	U	---
Alpha-chlordane	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---
Gamma-chlordane	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---
Toxaphene	330	U	---	320	U	---	320	U	---	340	U	---	320	U	---	380	U	---	350	U	---
Aroclor-1016	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---
Aroclor-1221	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---
Aroclor-1232	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---
Aroclor-1242	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA4-11A-SS	Q*	DV**	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**	SA4-12B-SS	Q*	DV**
Aroclor-1248	160	U	---	160	U	---	160	U	---	170	U	---	160	U	---	190	U	---	180	U	---
Aroclor-1254	330	U	---	320	U	---	320	U	---	340	U	---	320	U	---	380	U	---	350	U	---
Aroclor-1260	330	U	---	320	U	---	320	U	---	340	U	---	320	U	---	380	U	---	350	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**	SA4-13D-SS	Q*	DV**
Alpha-BHC	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Beta-BHC	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Delta-BHC	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Gamma-BHC	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Heptachlor	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Aldrin	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Heptachlor epoxide	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Endosulfan I	18	U	---	19	U	---	17	U	---	16	U	---	17	U	---	17	U	---	18	U	---
Dieldrin	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
4,4'-DDE	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
Endrin	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
Endosulfan II	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
4,4'-DDD	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
Endosulfan sulfate	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
4,4'-DDT	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
Methoxychlor	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---
Endrin ketone	37	U	---	37	U	---	34	U	---	33	U	---	34	U	---	33	U	---	36	U	---
Alpha-chlordane	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---
Gamma-chlordane	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---
Toxaphene	370	U	---	370	U	---	340	U	---	330	U	---	340	U	---	330	U	---	360	U	---
Aroclor-1016	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---
Aroclor-1221	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---
Aroclor-1232	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---
Aroclor-1242	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**	SA4-13D-SS	Q*	DV**
Aroclor-1248	180	U	---	190	U	---	170	U	---	160	U	---	170	U	---	170	U	---	180	U	---
Aroclor-1254	370	U	---	370	U	---	340	U	---	330	U	---	340	U	---	330	U	---	360	U	---
Aroclor-1260	370	U	---	370	U	---	340	U	---	330	U	---	340	U	---	330	U	---	360	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**	SA4-15A-SS	Q*	DV**
Alpha-BHC	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Beta-BHC	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Delta-BHC	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Gamma-BHC	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Heptachlor	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Aldrin	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Heptachlor epoxide	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Endosulfan I	19	U	---	17	U	---	18	U	---	17	U	---	17	U	---	17	U	---	17	U	---
Dieldrin	320		---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
4,4'-DDE	37	U	---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
Endrin	37	U	---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
Endosulfan II	37	U	---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
4,4'-DDD	37	U	---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
Endosulfan sulfate	37	U	---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
4,4'-DDT	37	U	---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
Methoxychlor	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---
Endrin ketone	37	U	---	34	U	---	35	U	---	35	U	---	33	U	---	34	U	---	33	U	---
Alpha-chlordane	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---
Gamma-chlordane	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---
Toxaphene	370	U	---	340	U	---	350	U	---	350	U	---	330	U	---	340	U	---	330	U	---
Aroclor-1016	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1221	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1232	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1242	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**	SA4-15A-SS	Q*	DV**
Aroclor-1248	190	U	---	170	U	---	180	U	---	170	U	---	170	U	---	170	U	---	170	U	---
Aroclor-1254	370	U	---	340	U	---	350	U	---	350	U	---	330	U	---	340	U	---	330	U	---
Aroclor-1260	370	U	---	340	U	---	350	U	---	350	U	---	330	U	---	340	U	---	330	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**	SA5-16B-SS	Q*	DV**	SA5-16C-SS	Q*	DV**
Alpha-BHC	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Beta-BHC	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Delta-BHC	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Gamma-BHC	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Heptachlor	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Aldrin	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Heptachlor epoxide	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Endosulfan I	19	U	---	18	U	---	19	U	---	19	U	---	18	U	---	17	U	---	17	U	---
Dieldrin	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	35	U	---	35	U	---
4,4'-DDE	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	94		---	35	U	---
Endrin	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	35	U	---	35	U	---
Endosulfan II	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	35	U	---	35	U	---
4,4'-DDD	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	35	U	---	35	U	---
Endosulfan sulfate	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	35	U	---	35	U	---
4,4'-DDT	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	90		---	35	U	---
Methoxychlor	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---
Endrin ketone	38	U	---	35	U	---	37	U	---	37	U	---	36	U	---	35	U	---	35	U	---
Alpha-chlordane	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---
Gamma-chlordane	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---
Toxaphene	380	U	---	350	U	---	370	U	---	370	U	---	360	U	---	350	U	---	350	U	---
Aroclor-1016	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---
Aroclor-1221	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---
Aroclor-1232	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---
Aroclor-1242	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**	SA5-16B-SS	Q*	DV**	SA5-16C-SS	Q*	DV**
Aroclor-1248	190	U	---	180	U	---	190	U	---	190	U	---	180	U	---	170	U	---	170	U	---
Aroclor-1254	380	U	---	350	U	---	370	U	---	370	U	---	360	U	---	350	U	---	350	U	---
Aroclor-1260	380	U	---	350	U	---	370	U	---	370	U	---	360	U	---	350	U	---	350	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**	SA5-17D-SS	Q*	DV**	SA5-17E-SS	Q*	DV**
Alpha-BHC	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Beta-BHC	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Delta-BHC	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Gamma-BHC	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Heptachlor	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Aldrin	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Heptachlor epoxide	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Endosulfan I	17	U	---	210	U	---	18	U	---	18	U	---	17	U	---	18	U	---	18	U	---
Dieldrin	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	
4,4'-DDE	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	---
Endrin	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	---
Endosulfan II	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	---
4,4'-DDD	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	---
Endosulfan sulfate	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	---
4,4'-DDT	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	---
Methoxychlor	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---
Endrin ketone	34	U	---	410	U	---	36	U	---	36	U	---	35	U	---	37	U	---	36	U	---
Alpha-chlordane	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---
Gamma-chlordane	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---
Toxaphene	340	U	---	4100	U	---	360	U	---	360	U	---	350	U	---	370	U	---	360	U	---
Aroclor-1016	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---
Aroclor-1221	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---
Aroclor-1232	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---
Aroclor-1242	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**	SA5-17D-SS	Q*	DV**	SA5-17E-SS	Q*	DV**
Aroclor-1248	170	U	---	2100	U	---	180	U	---	180	U	---	170	U	---	180	U	---	180	U	---
Aroclor-1254	340	U	---	4100	U	---	360	U	---	360	U	---	350	U	---	370	U	---	360	U	---
Aroclor-1260	340	U	---	4100	U	---	360	U	---	360	U	---	350	U	---	370	U	---	360	U	---

Table A-5.2 Pesticides/PCB's

Compound (ug/Kg)	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**	REFC-19A-SS	Q*	DV**	REFC-19B-SS	Q*	DV**
Alpha-BHC	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	---
Beta-BHC	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	---
Delta-BHC	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	---
Gamma-BHC	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	--
Heptachlor	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	---
Aldrin	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	---
Heptachlor epoxide	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	---
Endosulfan I	18	U	---	18	U	---	18	U	---	18	U	---	18	U	---	17	U	---	18	U	---
Dieldrin	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
4,4'-DDE	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
Endrin	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
Endosulfan II	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
4,4'-DDD	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
Endosulfan sulfate	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
4,4'-DDT	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
Methoxychlor	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---
Endrin ketone	36	U	---	35	U	---	35	U	---	35	U	---	35	U	---	34	U	---	37	U	---
Alpha-chlordane	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---
Gamma-chlordane	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---
Toxaphene	360	U	---	350	U	---	350	U	---	350	U	---	350	U	---	340	U	---	370	U	---
Aroclor-1016	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---
Aroclor-1221	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---
Aroclor-1232	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---
Aroclor-1242	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's																							
	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**	REFC-19A-SS	Q*	DV**	REFC-19B-SS	Q*	DV**			
Aroclor-1248	180	U	---	180	U	---	180	U	---	180	U	---	180	U	---	170	U	---	180	U	---			
Aroclor-1254	360	U	---	350	U	---	350	U	---	350	U	---	350	U	---	340	U	---	370	U	---			
Aroclor-1260	360	U	---	350	U	---	350	U	---	350	U	---	350	U	---	340	U	---	370	U	---			

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's								
	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**
Alpha-BHC	22	U	---	16	U	---	17	U	---
Beta-BHC	22	U	---	16	U	---	17	U	---
Delta-BHC	22	U	---	16	U	---	17	U	---
Gamma-BHC	22	U	---	16	U	---	17	U	---
Heptachlor	22	U	---	16	U	---	17	U	---
Aldrin	22	U	---	16	U	---	17	U	---
Heptachlor epoxide	22	U	---	16	U	---	17	U	---
Endosulfan I	22	U	---	16	U	---	17	U	---
Dieldrin	44	U	---	33	U	---	34	U	---
4,4'-DDE	44	U	---	33	U	---	34	U	---
Endrin	44	U	---	33	U	---	34	U	---
Endosulfan II	44	U	---	33	U	---	34	U	---
4,4'-DDD	44	U	---	33	U	---	34	U	---
Endosulfan sulfate	44	U	---	33	U	---	34	U	---
4,4'-DDT	44	U	---	33	U	---	34	U	---
Methoxychlor	220	U	---	160	U	---	170	U	---
Endrin ketone	44	U	---	33	U	---	34	U	---
Alpha-chlordane	220	U	---	160	U	---	170	U	---
Gamma-chlordane	220	U	---	160	U	---	170	U	---
Toxaphene	440	U	---	330	U	---	340	U	---
Aroclor-1016	220	U	---	160	U	---	170	U	---
Aroclor-1221	220	U	---	160	U	---	170	U	---
Aroclor-1232	220	U	---	160	U	---	170	U	---
Aroclor-1242	220	U	---	160	U	---	170	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.2 Pesticides/PCB's								
	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**
Aroclor-1248	220	u	---	160	u	---	170	u	---
Aroclor-1254	440	u	---	330	u	---	340	u	---
Aroclor-1260	440	u	---	330	u	---	340	u	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA1-1A-SS	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1E-SS	Q*	DV**	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**
Chloromethane	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Bromomethane	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Vinyl Chloride	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Chloroethane	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Methylene Chloride	6	U	---	12	B	U	13	B	U	17	B	U	13	B	U	23	B	U	19	B	U
Acetone	55	B	UJ	55	B	UJ	50	B	UJ	45	B	UJ	58	B	UJ	67	B	UJ	35	B	UJ
Carbon Disulfide	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
1,1-Dichloroethene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
1,1-Dichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
1,2-Dichloroethene (Total)	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Chloroform	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
1,2-Dichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
2-Butanone	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
1,1,1-Trichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Carbon Tetrachloride	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Vinyl Acetate	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Bromodichloromethane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
1,2-Dichloropropane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
cis-1,3-Dichloropropene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Trichloroethene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Dibromochloromethane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
1,1,2-Trichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Benzene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
trans-1,3-Dichloropropene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	SA1-1A-SS	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1E-SS	Q*	DV**	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**
	Table A-5.3 Volatiles																				
Bromoform	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
4-Methyl-2-Pentanone	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
2-Hexanone	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Tetrachloroethene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
1,1,2,2-Tetrachlorethane	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Toluene	1	J	---	3	J	---	4	J	---	5	U	---	2	J	---	5	U	---	6	U	---
Chlorobenzene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Ethylbenzene	6	U	---	2	J	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Styrene	6	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	---	6	U	---
Xylene (Total)	2	J	---	5	J	---	13		---	2	J	---	4	J	---	5	U	---	3	J	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**
Chloromethane	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
Bromomethane	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
Vinyl Chloride	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
Chloroethane	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
Methylene Chloride	17	B	U	19	B	U	18	B	U	5	BJ	UJ	7	B	UJ	11	B	U	11	B	UJ
Acetone	23	B	UJ	37	B	UJ	30	B	UJ	33	B	UJ	56	B	UJ	56	B	UJ	40	B	UJ
Carbon Disulfide	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
1,1-Dichloroethene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
1,1-Dichloroethane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
1,2-Dichloroethene (Total)	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Chloroform	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
1,2-Dichloroethane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
2-Butanone	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
1,1,1-Trichloroethane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Carbon Tetrachloride	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Vinyl Acetate	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
Bromodichloromethane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
1,2-Dichloropropane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
cis-1,3-Dichloropropene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Trichloroethene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Dibromochloromethane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
1,1,2-Trichloroethane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Benzene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
trans-1,3-Dichloropropene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**
Bromoform	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
4-Methyl-2-Pentanone	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
2-Hexanone	11	U	---	12	U	---	11	U	---	10	U	R	12	U	R	11	U	---	11	U	R
Tetrachloroethene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
1,1,2,2-Tetrachlorethane	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Toluene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Chlorobenzene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Ethylbenzene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Styrene	5	U	---	6	U	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R
Xylene (Total)	5	U	---	2	J	---	6	U	---	5	U	R	6	U	R	5	U	---	5	U	R

Table A-5.3 Volatiles

Compound (ug/Kg)	SA1-3E-SS	Q*	DV**	SA1-4A-SS	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**
Chloromethane	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	6	U	---
Bromomethane	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	6	U	---
Vinyl Chloride	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	6	U	---
Chloroethane	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	6	U	---
Methylene Chloride	8	B	UJ	6	U	UJ	5	BJ	UJ	5	U	---	6	U	---	6	U	---	6	U	---
Acetone	45	B	UJ	41	B	UJ	59	B	UJ	24	B	UJ	25	B	UJ	30	B	UJ	12	U	---
Carbon Disulfide	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	12	U	---
1,1-Dichloroethene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
1,1-Dichloroethane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
1,2-Dichloroethene (Total)	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Chloroform	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
1,2-Dichloroethane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
2-Butanone	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	12	U	R
1,1,1-Trichloroethane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Carbon Tetrachloride	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Vinyl Acetate	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	12	U	---
Bromodichloromethane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
1,2-Dichloropropane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
cis-1,3-Dichloropropene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Trichloroethene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Dibromochloromethane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
1,1,2-Trichloroethane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Benzene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
trans-1,3-Dichloropropene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---

* Data Qualifier. ** Data Validation.

Table A-5.3 Volatiles

Compound (ug/Kg)	SA1-3E-SS	Q*	DV**	SA1-4A-SS	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**
Bromoform	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
4-Methyl-2-Pentanone	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	12	U	---
2-Hexanone	14	U	R	12	U	UJ	12	U	R	11	U	---	12	U	---	12	U	---	12	U	---
Tetrachloroethene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
1,1,2,2-Tetrachlorethane	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Toluene	7	U	R	6	U	UJ	2	J	---	5	U	---	6	U	---	6	U	---	6	U	---
Chlorobenzene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Ethylbenzene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	6	U	---
Styrene	7	U	R	6	U	UJ	6	U	R	5	U	---	6	U	---	6	U	---	12	U	---
Xylene (Total)	7	U	R	6	U	UJ	6	U	R	2	J	---	6	U	---	6	U	---	12	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA2-5B-SS	Q*	DV**	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**
Chloromethane	6	U	---	6	U	---	5	U	---	6	U	---	12	U	---	11	U	---	10	U	---
Bromomethane	6	U	---	6	U	---	5	U	---	6	U	---	12	U	---	11	U	---	10	U	---
Vinyl Chloride	6	U	---	6	U	---	5	U	---	6	U	---	12	U	---	11	U	---	10	U	---
Chloroethane	6	U	---	6	U	---	5	U	---	6	U	---	12	U	---	11	U	---	10	U	---
Methylene Chloride	6	U	---	6	U	---	5	U	---	6	U	---	5	BJ	U	4	BJ	5U	4	BJ	5U
Acetone	12	U	---	12	U	---	11	U	---	12	U	---	54	B	UJ	64	B	UJ	45	B	UJ
Carbon Disulfide	12	U	---	12	U	---	11	U	---	12	U	---	6	U	---	5	U	---	5	U	---
1,1-Dichloroethene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
1,1-Dichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
1,2-Dichloroethene (Total)	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
Chloroform	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U		5	U	---
1,2-Dichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
2-Butanone	12	U	R	12	U	R	11	U	R	12	U	R	12	U	---	11	U	---	10	U	---
1,1,1-Trichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
Carbon Tetrachloride	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
Vinyl Acetate	12	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	10	U	---
Bromodichloromethane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
1,2-Dichloropropane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
cis-1,3-Dichloropropene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
Trichloroethene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
Dibromochloromethane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
1,1,2-Trichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
Benzene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
trans-1,3-Dichloropropene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA2-5B-SS	Q*	DV**	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**
Bromoform	6	U	---	6	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---
4-Methyl-2-Pentanone	12	U	---	12	U	---	11	U	---	12	U	---	12	U	UJ	11	U	---	10	U	---
2-Hexanone	12	U	---	12	U	---	11	U	---	12	U	---	12	U	UJ	11	U	---	10	U	---
Tetrachloroethene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---
1,1,2,2-Tetrachlorethane	6	U	---	6	U	---	5	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---
Toluene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	UJ	5	U	---	2	J	---
Chlorobenzene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---
Ethylbenzene	6	U	---	6	U	---	5	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---
Styrene	12	U	---	12	U	---	11	U	---	12	U	---	6	U	UJ	5	U	---	5	U	---
Xylene (Total)	12	U	---	12	U	---	11	U	---	12	U	---	6	U	UJ	5	U	---	5	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA2-6D-SS	Q*	DV**	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**
Chloromethane	11	U	---	10	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Bromomethane	11	U	---	10	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Vinyl Chloride	11	U	---	10	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Chloroethane	11	U	---	10	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Methylene Chloride	4	BJ	5U	2	BJ	5U	6	U	---	6	U	---	1	J	---	5	U	---	5	U	---
Acetone	46	B	UJ	40	B	UJ	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---
Carbon Disulfide	5	U	---	5	U	---	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---
1,1-Dichloroethene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U		5	U	---
1,2-Dichloroethene (Total)	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Chloroform	5	U	---	5	U	---	5	J	6U	1	J	6U	7		---	5	U	---	2	J	5U
1,2-Dichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
2-Butanone	11	U	---	10	U	---	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---
1,1,1-Trichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Carbon Tetrachloride	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Vinyl Acetate	11	U	---	10	U	---	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---
Bromodichloromethane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloropropane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
cis-1,3-Dichloropropene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Trichloroethene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Dibromochloromethane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,1,2-Trichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Benzene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
trans-1,3-Dichloropropene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA2-6D-SS	Q*	DV**	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**
Bromoform	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
4-Methyl-2-Pentanone	11	U	---	10	U	---	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---
2-Hexanone	11	U	---	10	U	---	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---
Tetrachloroethene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,1,2,2-Tetrachlorethane	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Toluene	2	J	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Chlorobenzene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Ethylbenzene	5	U	---	5	U	---	6	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Styrene	5	U	---	5	U	---	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---
Xylene (Total)	5	U	---	5	U	---	12	U	---	12	U	---	11	U	---	11	U	---	10	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA3-8A-SS	Q*	DV**	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**
Chloromethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Bromomethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Vinyl Chloride	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Chloroethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Methylene Chloride	6	U	---	2	J	---	3	J	---	5	U	---	6	U	---	6	U	---	6	U	---
Acetone	12	U	---	11	U	---	10	U	---	11	U	---	11	U	---	12	U	---	12	U	---
Carbon Disulfide	12	U	---	11	U	---	10	U	---	11	U	---	11	U	---	12	U	---	12	U	---
1,1-Dichloroethene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
1,1-Dichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
1,2-Dichloroethene (Total)	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Chloroform	6	U	---	5	U	---	5	U	---	3	J	5U	6	U	---	6	U	---	6	U	---
1,2-Dichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
2-Butanone	12	U	---	11	U	---	10	U	---	11	U	R	11	U	---	12	U	R	12	U	R
1,1,1-Trichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Carbon Tetrachloride	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Vinyl Acetate	12	U	---	11	U	---	10	U	---	11	U	---	11	U	---	12	U	---	12	U	---
Bromodichloromethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
1,2-Dichloropropane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
cis-1,3-Dichloropropene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Trichloroethene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Dibromochloromethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
1,1,2-Trichloroethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Benzene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
trans-1,3-Dichloropropene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA3-8A-SS	Q*	DV**	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**
Bromoform	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
4-Methyl-2-Pentanone	12	U	---	11	U	---	10	U	---	11	U	---	11	U	---	12	U	---	12	U	---
2-Hexanone	12	U	---	11	U	---	10	U	---	11	U	---	11	U	---	12	U	---	12	U	---
Tetrachloroethene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
1,1,2,2-Tetrachlorethane	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Toluene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Chlorobenzene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Ethylbenzene	6	U	---	5	U	---	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---
Styrene	12	U	---	11	U	---	10	U	---	11	U	---	11	U	---	12	U	---	12	U	---
Xylene (Total)	12	U	---	11	U	---	10	U	---	11	U	---	11	U	---	12	U	---	12	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA3-9C-SS	Q*	DV**	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**
Chloromethane	6	U	---	6	U	---	6	U	---	11	U	---	12	U	---	11	U	---	11	U	UJ
Bromomethane	6	U	---	6	U	---	6	U	---	11	U	---	12	U	---	11	U	---	11	U	UJ
Vinyl Chloride	6	U	---	6	U	---	6	U	---	11	U	---	12	U	---	11	U	---	11	U	UJ
Chloroethane	6	U	---	6	U	---	6	U	---	11	U	---	12	U	---	11	U	---	11	U	UJ
Methylene Chloride	6	U	---	6	U	---	6	U	---	24	B	U	6	U	---	32	B	U	30	B	UJ
Acetone	11	U	---	12	U	---	13	U	---	100	B	UJ	110	B	UJ	120	B	UJ	180	B	UJ
Carbon Disulfide	11	U	---	12	U	---	13	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
1,1-Dichloroethene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
1,1-Dichloroethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
1,2-Dichloroethene (Total)	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Chloroform	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
1,2-Dichloroethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
2-Butanone	11	U	R	12	U	R	13	U	R	11	U	---	12	U	---	11	U	---	11	U	UJ
1,1,1-Trichloroethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Carbon Tetrachloride	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Vinyl Acetate	11	U	---	12	U	---	13	U	---	11	U	---	12	U	---	11	U	---	11	U	UJ
Bromodichloromethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
1,2-Dichloropropane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
cis-1,3-Dichloropropene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Trichloroethene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Dibromochloromethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
1,1,2-Trichloroethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Benzene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA3-9C-SS	Q*	DV**	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**
trans-1,3-Dichloropropene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Bromoform	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
4-Methyl-2-Pentanone	11	U	---	12	U	---	13	U	---	11	U	---	12	U	---	11	U	---	11	U	UJ
2-Hexanone	11	U	---	12	U	---	13	U	---	11	U	---	12	U	---	11	U	---	11	U	UJ
Tetrachloroethene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
1,1,2,2-Tetrachlorethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Toluene	6	U	---	6	U	---	6	U	---	2	J	---	2	J	---	6	U	---	2	J	---
Chlorobenzene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Ethylbenzene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Styrene	11	U	---	12	U	---	13	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ
Xylene (Total)	11	U	---	12	U	---	13	U	---	6	U	---	6	U	---	6	U	---	6	U	UJ

Table A-5.3 Volatiles

Compound (ug/Kg)	SA3-10E-SS	Q*	DV**	SA4-11A-SS	Q*	DV**	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**
Chloromethane	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
Bromomethane	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
Vinyl Chloride	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
Chloroethane	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
Methylene Chloride	30	B	U	5	BJ	UJ	6	B	UJ	7	B	UJ	5	U	R	9	B	UJ	7	B	U
Acetone	110	B	UJ	36	B	UJ	44	B	UJ	58	B	UJ	49	B	UJ	33	B	UJ	38	B	UJ
Carbon Disulfide	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
1,1-Dichloroethene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
1,1-Dichloroethane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
1,2-Dichloroethene (Total)	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Chloroform	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
1,2-Dichloroethane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
2-Butanone	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
1,1,1-Trichloroethane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Carbon Tetrachloride	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Vinyl Acetate	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
Bromodichloromethane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
1,2-Dichloropropane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
cis-1,3-Dichloropropene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Trichloroethene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Dibromochloromethane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
1,1,2-Trichloroethane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Benzene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---

* Data Qualifier. ** Data Validation.

Table A-5.3 Volatiles

Compound (ug/Kg)	SA3-10E-SS	Q*	DV**	SA4-11A-SS	Q*	DV**	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**
trans-1,3-Dichloropropene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Bromoform	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
4-Methyl-2-Pentanone	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
2-Hexanone	12	U	---	10	U	R	10	U	R	10	U	R	11	U	R	10	U	R	12	U	---
Tetrachloroethene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
1,1,2,2-Tetrachlorethane	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Toluene	6	U	---	3	J	---	1	J	---	5	U	R	9		J	2	J	---	6	U	---
Chlorobenzene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Ethylbenzene	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---
Styrene	6	U	---	5	U	R	5	U	R	5	U	R	25		J	5	U	R	6	U	---
Xylene (Total)	6	U	---	5	U	R	5	U	R	5	U	R	5	U	R	5	U	R	6	U	---

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA4-12B-SS	Q*	DV**	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**
Chloromethane	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	---	11	U	---	10	U	---
Bromomethane	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	---	11	U	---	10	U	---
Vinyl Chloride	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	---	11	U	---	10	U	---
Chloroethane	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	---	11	U	---	10	U	---
Methylene Chloride	5	BJ	6U	4	BJ	6U	6	B	UJ	5	BJ	U	5	U	---	4	BJ	5U	5	U	---
Acetone	29	B	UJ	32	B	U	63	B	UJ	47	B	UJ	31	B	UJ	30	B	UJ	10	U	UJ
Carbon Disulfide	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethene (Total)	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Chloroform	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
2-Butanone	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	UJ	11	U	UJ	10	U	UJ
1,1,1-Trichloroethane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Carbon Tetrachloride	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Vinyl Acetate	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	---	11	U	---	10	U	---
Bromodichloromethane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloropropane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
cis-1,3-Dichloropropene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Trichloroethene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Dibromochloromethane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2-Trichloroethane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Benzene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA4-12B-SS	Q*	DV**	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**
trans-1,3-Dichloropropene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Bromoform	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
4-Methyl-2-Pentanone	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	---	11	U	---	10	U	---
2-Hexanone	11	U	---	11	U	---	12	U	UJ	11	U	---	10	U	---	11	U	---	10	U	---
Tetrachloroethene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
1,1,2,2-Tetrachlorethane	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Toluene	6	U	---	6	U	---	6	U	UJ	5	U	---	2	J	---	5	U	---	5	U	---
Chlorobenzene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Ethylbenzene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Styrene	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---
Xylene (Total)	6	U	---	6	U	---	6	U	UJ	5	U	---	5	U	---	5	U	---	5	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA4-13D-SS	Q*	DV**	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**
Chloromethane	11	U	---	12	U	---	11	U	---	11	U	---	11	U	---	10	U	---	11	U	---
Bromomethane	11	U	---	12	U	---	11	U	---	11	U	---	11	U	---	10	U		11	U	---
Vinyl Chloride	11	U	---	12	U	---	11	U	---	11	U	---	11	U	---	10	U	---	11	U	---
Chloroethane	11	U	---	12	U	---	11	U	---	11	U	---	11	U	---	10	U	---	11	U	---
Methylene Chloride	5	BJ	U	5	BJ	U	8		U	8		U	11		U	4	BJ	5U	4	BJ	5U
Acetone	29	B	UJ	56	B	UJ	19	B	UJ	74	B	UJ	71	B	UJ	45	B	UJ	40	B	UJ
Carbon Disulfide	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,1-Dichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethene (Total)	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Chloroform	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
2-Butanone	11	U	UJ	12	U	UJ	11	U	---	11	U	---	11	U	---	10	U	---	11	U	---
1,1,1-Trichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Carbon Tetrachloride	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Vinyl Acetate	11	U	---	12	U	---	11	U	---	11	U	---	11	U	---	10	U	---	11	U	---
Bromodichloromethane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,2-Dichloropropane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
cis-1,3-Dichloropropene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Trichloroethene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Dibromochloromethane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,1,2-Trichloroethane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Benzene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA4-13D-SS	Q*	DV**	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**
trans-1,3-Dichloropropene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Bromoform	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
4-Methyl-2-Pentanone	11	U	---	12	U	---	11	U	---	11	U	---	11	U	---	10	U	---	11	U	---
2-Hexanone	11	U	---	12	U	---	11	U	---	11	U	---	11	U	---	10	U	---	11	U	---
Tetrachloroethene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
1,1,2,2-Tetrachlorethane	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Toluene	6	U	---	6	U	---	1	J	---	6	U	---	5	U	---	2	J	---	5	U	---
Chlorobenzene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Ethylbenzene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Styrene	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---
Xylene (Total)	6	U	---	6	U	---	5	U	---	6	U	---	5	U	---	5	U	---	5	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA4-15A-SS	Q*	DV**	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**	SA5-16B-SS	Q*	DV**
Chloromethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Bromomethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Vinyl Chloride	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Chloroethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Methylene Chloride	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Acetone	10	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	11	U	---
Carbon Disulfide	10	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	11	U	---
1,1-Dichloroethene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,1-Dichloroethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,2-Dichloroethene (Total)	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Chloroform	5	U	---	6	U	---	5	U	---	9		U	6	U	---	6	U	---	5	U	---
1,2-Dichloroethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
2-Butanone	10	U	R	12	U	R	11	U	R	12	U	R	12	U	R	11	U	R	11	U	R
1,1,1-Trichloroethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Carbon Tetrachloride	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Vinyl Acetate	10	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	11	U	---
Bromodichloromethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,2-Dichloropropane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
cis-1,3-Dichloropropene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Trichloroethene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Dibromochloromethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,1,2-Trichloroethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Benzene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---

* Data Qualifier. ** Data Validation.

Table A-5.3 Volatiles

Compound (ug/Kg)	SA4-15A-SS	Q*	DV**	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**	SA5-16B-SS	Q*	DV**
trans-1,3-Dichloropropene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Bromoform	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
4-Methyl-2-Pentanone	10	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	11	U	---
2-Hexanone	10	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	11	U	---
Tetrachloroethene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,1,2,2-Tetrachlorethane	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Toluene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Chlorobenzene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Ethylbenzene	5	U	---	6	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Styrene	10	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	11	U	---
Xylene (Total)	10	U	---	12	U	---	11	U	---	12	U	---	12	U	---	11	U	---	11	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	SA5-16C-SS	Q*	DV**	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**	SA5-17D-SS	Q*	DV**
Chloromethane	5	U	---	5	U	---	6	U	---	11	U	---	11	U	---	11	U	---	12	U	UJ
Bromomethane	5	U	---	5	U	---	6	U	---	11	U	---	11	U	---	11	U	---	12	U	UJ
Vinyl Chloride	5	U	---	5	U	---	6	U	---	11	U	---	11	U	---	11	U	---	12	U	UJ
Chloroethane	5	U	---	5	U	---	6	U	---	11	U	---	11	U	---	11	U	---	12	U	UJ
Methylene Chloride	5	U	---	5	U	---	6	U	---	7	B	U	7	B	U	7	B	U	9	B	U
Acetone	11	U	---	11	U	---	13	U	---	45	B	UJ	58	B	UJ	73	B	UJ	50	B	J
Carbon Disulfide	11	U	---	11	U	---	13	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
1,1-Dichloroethene	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
1,1-Dichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
1,2-Dichloroethene (Total)	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
Chloroform	6		---	5	J	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
1,2-Dichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
2-Butanone	11	U	R	11	U	R	13	U	---	11	U	---	11	U	---	11	U	---	12	U	UJ
1,1,1-Trichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
Carbon Tetrachloride	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
Vinyl Acetate	11	U	---	11	U	---	13	U	---	11	U	---	11	U	---	11	U	---	12	U	UJ
Bromodichloromethane	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
1,2-Dichloropropane	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
cis-1,3-Dichloropropene	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
Trichloroethene	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
Dibromochloromethane	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
1,1,2-Trichloroethane	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
Benzene	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.3 Volatiles																							
	SA5-16C-SS	Q*	DV**	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**	SA5-17D-SS	Q*	DV**
trans-1,3-Dichloropropene	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
Bromoform	5	U	---	5	U	---	6	U	---	6	U	---	6	U	---	5	U	---	6	U	UJ
4-Methyl-2-Pentanone	11	U	---	11	U	---	13	U	---	11	U	---	11	U	---	11	U	---	12	U	UJ
2-Hexanone	11	U	---	11	U	---	13	U	---	11	U	UJ	11	U	---	11	U	---	12	U	UJ
Tetrachloroethene	5	U	---	5	U	---	6	U	---	6	U	UJ	6	U	---	5	U	---	6	U	UJ
1,1,2,2-Tetrachlorethane	5	U	---	5	U	---	6	U	---	6	U	UJ	6	U	---	5	U	---	6	U	UJ
Toluene	5	U	---	5	U	---	6	U	---	6	U	UJ	6	U	---	5	U	---	2	J	---
Chlorobenzene	5	U	---	5	U	---	6	U	---	6	U	UJ	6	U	---	5	U	---	6	U	UJ
Ethylbenzene	5	U	---	5	U	---	6	U	---	6	U	UJ	6	U	---	5	U	---	6	U	UJ
Styrene	11	U	---	11	U	---	13	U	---	6	U	UJ	6	U	---	5	U	---	6	U	UJ
Xylene (Total)	11	U	---	11	U	---	13	U	---	6	U	UJ	6	U	---	5	U	---	6	U	UJ

Table A-5.3 Volatiles

Compound (ug/Kg)	SA5-17E-SS	Q*	DV**	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**	REFC-19A-SS	Q*	DV**
Chloromethane	11	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Bromomethane	11	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Vinyl Chloride	11	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Chloroethane	11	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Methylene Chloride	8	B	U	6	U	---	6	U	---	6	U	--	6	U	---	6	U	---	5	U	---
Acetone	73	B	UJ	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Carbon Disulfide	6	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
1,1-Dichloroethene	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,1-Dichloroethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,2-Dichloroethene (Total)	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Chloroform	6	U	---	6	U	---	7		---	6		---	6	U	---	4	J	---	5	U	---
1,2-Dichloroethane	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
2-Butanone	11	U	---	11	U	R	11	U	R	11	U	R	11	U	R	11	U	R	11	U	---
1,1,1-Trichloroethane	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Carbon Tetrachloride	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Vinyl Acetate	11	U	UJ	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Bromodichloromethane	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,2-Dichloropropane	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
cis-1,3-Dichloropropene	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Trichloroethene	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Dibromochloromethane	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,1,2-Trichloroethane	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Benzene	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---

* Data Qualifier. ** Data Validation.

Table A-5.3 Volatiles

Compound (ug/Kg)	SA5-17E-SS	Q*	DV**	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**	REFC-19A-SS	Q*	DV**
trans-1,3-Dichloropropene	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Bromoform	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
4-Methyl-2-Pentanone	11	U	UJ	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
2-Hexanone	11	U	UJ	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Tetrachloroethene	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
1,1,2,2-Tetrachlorethane	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Toluene	2	J	---	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Chlorobenzene	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Ethylbenzene	6	U	UJ	6	U	---	6	U	---	6	U	---	6	U	---	6	U	---	5	U	---
Styrene	6	U	UJ	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---
Xylene (Total)	6	U	UJ	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---	11	U	---

Table A-5.3 Volatiles

Compound (ug/Kg)	REFC-19B-SS	Q*	DV**	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**	TB4/30/93	Q*	DV**	TRIP BLANK (2)	Q*	DV**	TRIP BLANK (6)	Q*	DV**
Chloromethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	10	U	---	10	U	R
Bromomethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	10	U	---	10	U	R
Vinyl Chloride	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	10	U	---	10	U	R
Chloroethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	10	U	---	10	U	R
Methylene Chloride	6	U	---	16		---	9		---	5	U	---	5	U	---	15	B	U	7	B	UJ
Acetone	12	U	---	13	U	---	11	U	---	11	U	---	10	U	---	23	B	UJ	10	U	R
Carbon Disulfide	12	U	---	13	U	---	11	U	---	11	U	---	10	U	---	5	U	---	5	U	R
1,1-Dichloroethene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
1,1-Dichloroethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
1,2-Dichloroethene (Total)	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Chloroform	6	U	---	7	U	---	5	U	---	5	U	---	2	BJ	5U	5	U	---	5	U	R
1,2-Dichloroethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
2-Butanone	12	U	---	13	U	---	11	U	---	11	U	R	10	U	R	10	U	R	10	U	R
1,1,1-Trichloroethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Carbon Tetrachloride	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Vinyl Acetate	12	U	---	13	U	---	11	U	---	11	U	---	10	U	---	10	U	---	10	U	R
Bromodichloromethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U		5	U	R
1,2-Dichloropropane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
cis-1,3-Dichloropropene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Trichloroethene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Dibromochloromethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
1,1,2-Trichloroethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U		5	U	R
Benzene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R

Table A-5.3 Volatiles

Compound (ug/Kg)	REFC-19B-SS	Q*	DV**	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**	TB4/30/93	Q*	DV**	TRIP BLANK (2)	Q*	DV**	TRIP BLANK (6)	Q*	DV**
trans-1,3-Dichloropropene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Bromoform	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
4-Methyl-2-Pentanone	12	U	---	13	U	---	11	U	---	11	U	---	10	U	---	10	U	---	10	U	R
2-Hexanone	12	U	---	13	U	---	11	U	---	11	U	---	10	U	---	10	U	---	10	U	R
Tetrachloroethene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
1,1,2,2-Tetrachlorethane	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Toluene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Chlorobenzene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Ethylbenzene	6	U	---	7	U	---	5	U	---	5	U	---	5	U	---	5	U	---	5	U	R
Styrene	12	U	---	13	U	---	11	U	---	11	U	---	10	U	---	5	U	---	5	U	R
Xylene (Total)	12	U	---	13	U	---	11	U	---	11	U	---	10	U	---	5	U	---	5	U	R

Compound (ug/Kg)	Table A-5.3 Volatiles					
	TRIP BLANK (6)	Q*	DV**	TRIP BLANK (3)	Q*	DV**
Chloromethane	10	U	---	10	U	NP
Bromomethane	10	U	---	10	U	NP
Vinyl Chloride	10	U	---	10	U	NP
Chloroethane	10	U	---	10	U	NP
Methylene Chloride	18	B	U	7	B	NP
Acetone	25	B	UJ	10	U	NP
Carbon Disulfide	5	U	---	5	U	NP
1,1-Dichloroethene	5	U	---	5	U	NP
1,1-Dichloroethane	5	U	---	5	U	NP
1,2-Dichloroethene (Total)	5	U	---	5	U	NP
Chloroform	5	U	---	5	U	NP
1,2-Dichloroethane	5	U	---	5	U	NP
2-Butanone	10	U	---	10	U	NP
1,1,1-Trichloroethane	5	U	---	5	U	NP
Carbon Tetrachloride	5	U	---	5	U	NP
Vinyl Acetate	10	U	---	10	U	NP
Bromodichloromethane	5	U	---	5	U	NP
1,2-Dichloropropane	5	U	---	5	U	NP
cis-1,3-Dichloropropene	5	U	---	5	U	NP
Trichloroethene	5	U	---	5	U	NP
Dibromochloromethane	5	U	---	5	U	NP
1,1,2-Trichloroethane	5	U	---	5	U	NP
Benzene	5	U	---	5	U	NP

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.3 Volatiles					
	TRIP BLANK (6)	Q*	DV**	TRIP BLANK (3)	Q*	DV**
trans-1,3-Dichloropropene	5	U	---	5	U	NP
Bromoform	5	U	---	5	U	NP
4-Methyl-2-Pentanone	10	U	---	10	U	NP
2-Hexanone	10	U	---	10	U	NP
Tetrachloroethene	5	U	---	5	U	NP
1,1,2,2-Tetrachlorethane	5	U	---	5	U	NP
Toluene	5	U	---	5	U	NP
Chlorobenzene	5	U	---	5	U	NP
Ethylbenzene	5	U	---	5	U	NP
Styrene	5	U	---	5	U	NP
Xylene (Total)	5	U	---	5	U	NP

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA1-1A-SS	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1E-SS	Q*	DV**	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**
Phenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
bis (2-Cloroethyl)Ether	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2-Chlorophenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
1,3-Dichlorobenzene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
1,4-Dichlorobenzene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Benzyl Alcohol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
1,2-Dichlorobenzene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2-Methylphenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
bis (2-Chloroisopropyl)Ether	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
4-Methylphenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
N-Nitroso-Di-n-Propylamine	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Hexachloroethane	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Nitrobenzene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Isophorone	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2-Nitrophenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2,4-Dimethylphenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Benzoic Acid	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
bis (2-Chloroethoxy)Methane	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2,4-Dichlorophenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
1,2,4-Trichlorobenzene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Naphthalene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
4-Chloroaniline	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Hexachlorobutadiene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA1-1A-SS	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1E-SS	Q*	DV**	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**
4-Chloro-3-Methylphenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2-Methylnaphthalene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Hexachlorocyclopentadiene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2,4,6-Trichlorophenol	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2,4,5-Trichlorophenol	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
2-Chloronaphthalene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2-Nitroaniline	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
Dimethyl Phthalate	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Acenaphthylene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2,6-Dinitrotoluene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
3-Nitroaniline	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
Acenaphthene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2,4-Dinitrophenol	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
4-Nitrophenol	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
Dibenzofuran	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
2,4-Dinitrotoluene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Diethylphthalate	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
4-Chlorophenyl-phenylether	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Fluorene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
4-Nitroaniline	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
4,6-Dinitro-2-Methylphenol	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
N-Nitrosodiphenylamine (1)	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
4-Bromophenyl-phenylether	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA1-1A-SS	Q*	DV**	SA1-1B-SS	Q*	DV**	SA1-1C-SS	Q*	DV**	SA1-1D-SS	Q*	DV**	SA1-1E-SS	Q*	DV**	SA1-2A-SS	Q*	DV**	SA1-2B-SS	Q*	DV**
Hexachlorobenzene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Pentachlorophenol	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---	1800	U	---	1800	U	NP
Phenanthrene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Anthracene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Di-n-Butylphthalate	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Fluoranthrene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Pyrene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Butylbenzylphthalate	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
3,3'-Dichlorbenzidine	750	U	---	710	U	---	690	U	---	720	U	---	690	U	---	720	U	---	760	U	NP
Benzo(a)Anthracene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Chrysene	370	U	---	350	U	---	350	U	---	360	U	---	340	U	---	360	U	---	380	U	NP
Bis (2-Ethylhexyl)Phthalate	1700		---	730		---	300	J	---	1300		---	1200		---	180	J	---	470		NP
Di-n-Octyl Phthalate	370	U	---	350	U	---	350	U	UJ	360	U	---	340	U	---	360	U	---	380	U	NP
Benzo(b)Fluoranthene	370	U	---	350	U	---	350	U	UJ	360	U	---	340	U	---	360	U	---	380	U	NP
Benzo(k)Fluoranthene	370	U	---	350	U	---	350	U	UJ	360	U	---	340	U	---	360	U	---	380	U	NP
Benzo(a)Pyrene	370	U	---	350	U	---	350	U	UJ	360	U	---	340	U	---	360	U	---	380	U	NP
Indeno(1,2,3-cd)Pyrene	370	U	---	350	U	---	350	U	UJ	360	U	---	340	U	---	360	U	---	380	U	NP
Dibenz(a,h)Anthracene	370	U	---	350	U	---	350	U	UJ	360	U	---	340	U	---	360	U	---	380	U	NP
Benzo(g,h,i)Perylene	370	U	UJ	350	U	---	350	U	UJ	360	U	---	340	U	---	360	U	---	380	U	NP

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**
Phenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
bis (2-Cloroethyl)Ether	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2-Chlorophenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
1,3-Dichlorobenzene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
1,4-Dichlorobenzene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Benzyl Alcohol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	J
1,2-Dichlorobenzene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2-Methylphenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
bis (2-Chloroisopropyl)Ether	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
4-Methylphenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
N-Nitroso-Di-n-Propylamine	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Hexachloroethane	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Nitrobenzene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Isophorone	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2-Nitrophenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2,4-Dimethylphenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Benzoic Acid	1700	U	---	71	J	---	61	J	---	240	J	---	1700	U	---	1700	U	---	1700	U	---
bis (2-Chloroethoxy)Methane	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2,4-Dichlorophenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
1,2,4-Trichlorobenzene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Naphthalene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
4-Chloroaniline	350	U	UJ	390	U	UJ	380	U	UJ	420	U	---	350	U	UJ	350	U	UJ	360	U	---
Hexachlorobutadiene	350	U	---	390	U	---	380	U	---	420	U	UJ	350	U	---	350	U	---	360	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**
4-Chloro-3-Methylphenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2-Methylnaphthalene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Hexachlorocyclopentadiene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2,4,6-Trichlorophenol	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2,4,5-Trichlorophenol	1700	U	---	1900	U	---	1800	U	---	2100	U	---	1700	U	---	1700	U	---	1700	U	---
2-Chloronaphthalene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2-Nitroaniline	1700	U	---	1900	U	---	1800	U	---	2100	U	---	1700	U	---	1700	U	---	1700	U	---
Dimethyl Phthalate	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Acenaphthylene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2,6-Dinitrotoluene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
3-Nitroaniline	1700	U	UJ	1900	U	UJ	1800	U	UJ	2100	U	UJ	1700	U	UJ	1700	U	UJ	1700	U	---
Acenaphthene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2,4-Dinitrophenol	1700	U	---	1900	U	---	1800	U	---	2100	U	---	1700	U	---	1700	U	---	1700	U	---
4-Nitrophenol	1700	U	---	1900	U	---	1800	U	---	2100	U	---	1700	U	---	1700	U	---	1700	U	---
Dibenzofuran	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
2,4-Dinitrotoluene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Diethylphthalate	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
4-Chlorophenyl-phenylether	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Fluorene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
4-Nitroaniline	1700	U	---	1900	U	---	1800	U	---	2100	U	---	1700	U	---	1700	U	---	1700	U	---
4,6-Dinitro-2-Methylphenol	1700	U	---	1900	U	---	1800	U	---	2100	U	---	1700	U	---	1700	U	---	1700	U	---
N-Nitrosodiphenylamine (1)	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
4-Bromophenyl-phenylether	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA1-2C-SS	Q*	DV**	SA1-2D-SS	Q*	DV**	SA1-2E-SS	Q*	DV**	SA1-3A-SS	Q*	DV**	SA1-3B-SS	Q*	DV**	SA1-3C-SS	Q*	DV**	SA1-3D-SS	Q*	DV**
Hexachlorobenzene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Pentachlorophenol	1700	U	---	1900	U	---	1800	U	---	2100	U	---	1700	U	---	1700	U	---	1700	U	---
Phenanthrene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Anthracene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Di-n-Butylphthalate	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Fluoranthrene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Pyrene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Butylbenzylphthalate	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
3,3'-Dichlorbenzidine	710	U	---	780	U	---	760	U	---	850	U	---	710	U	---	690	U	---	710	U	---
Benzo(a)Anthracene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Chrysene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Bis (2-Ethylhexyl)Phthalate	200	J	---	480		---	360	J	---	140	J	---	60	J	---	150	J	---	150	J	---
Di-n-Octyl Phthalate	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Benzo(b)Fluoranthene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Benzo(k)Fluoranthene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Benzo(a)Pyrene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Indeno(1,2,3-cd)Pyrene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Dibenz(a,h)Anthracene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---
Benzo(g,h,i)Perylene	350	U	---	390	U	---	380	U	---	420	U	---	350	U	---	350	U	---	360	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA1-4A-SS	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**	SA2-5B-SS	Q*	DV**
Phenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
bis (2-Cloroethyl)Ether	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
2-Chlorophenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
1,3-Dichlorobenzene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
1,4-Dichlorobenzene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Benzyl Alcohol	410	U	---	380	U	---	350	U	---	400	U	J	410	U	---	390	U	UJ	380	U	UJ
1,2-Dichlorobenzene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
2-Methylphenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
bis (2-Chloroisopropyl)Ether	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	UJ
4-Methylphenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
N-Nitroso-Di-n-Propylamine	410	U	UJ	380	U	UJ	350	U	UJ	400	U	---	410	U	UJ	390	U	---	380	U	---
Hexachloroethane	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Nitrobenzene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Isophorone	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
2-Nitrophenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
2,4-Dimethylphenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Benzoic Acid	110	J	---	1800	U	---	1700	U	---	230	J	---	170	J	---	1900	U	---	87	J	---
bis (2-Chloroethoxy)Methane	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
2,4-Dichlorophenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
1,2,4-Trichlorobenzene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Naphthalene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
4-Chloroaniline	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Hexachlorobutadiene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA1-4A-SS	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**	SA2-5B-SS	Q*	DV**			
4-Chloro-3-Methylphenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
2-Methylnaphthalene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
Hexachlorocyclopentadiene	410	U	R	380	U	R	350	U	R	400	U	---	410	U	R	390	U	---	380	U	---			
2,4,6-Trichlorophenol	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
2,4,5-Trichlorophenol	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1900	U	---			
2-Chloronaphthalene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
2-Nitroaniline	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1900	U	---			
Dimethyl Phthalate	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
Acenaphthylene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
2,6-Dinitrotoluene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
3-Nitroaniline	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1900	U	---			
Acenaphthene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
2,4-Dinitrophenol	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	UJ	1900	U	---			
4-Nitrophenol	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1900	U	---			
Dibenzofuran	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
2,4-Dinitrotoluene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
Diethylphthalate	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
4-Chlorophenyl-phenylether	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
Fluorene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
4-Nitroaniline	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1900	U	---			
4,6-Dinitro-2-Methylphenol	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1900	U	---			
N-Nitrosodiphenylamine (1)	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			
4-Bromophenyl-phenylether	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---			

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA1-4A-SS	Q*	DV**	SA1-4B-SS	Q*	DV**	SA1-4C-SS	Q*	DV**	SA1-4D-SS	Q*	DV**	SA1-4E-SS	Q*	DV**	SA2-5A-SS	Q*	DV**	SA2-5B-SS	Q*	DV**
Hexachlorobenzene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Pentachlorophenol	2000	U	---	1800	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1900	U	---
Phenanthrene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Anthracene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Di-n-Butylphthalate	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Fluoranthrene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Pyrene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Butylbenzylphthalate	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
3,3'-Dichlorbenzidine	820	U	---	760	U	---	710	U	---	790	U	---	820	U	---	780	U	---	760	U	---
Benzo(a)Anthracene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Chrysene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Bis (2-Ethylhexyl)Phthalate	44	J	---	77	J	---	170	J	---	290	J	---	220	J	---	63	J	---	54	J	---
Di-n-Octyl Phthalate	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Benzo(b)Fluoranthene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Benzo(k)Fluoranthene	410	U	UJ	380	U	UJ	350	U	UJ	400	U	---	410	U	UJ	390	U	---	380	U	---
Benzo(a)Pyrene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Indeno(1,2,3-cd)Pyrene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Dibenz(a,h)Anthracene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---
Benzo(g,h,i)Perylene	410	U	---	380	U	---	350	U	---	400	U	---	410	U	---	390	U	---	380	U	---

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**	SA2-6D-SS	Q*	DV**
Phenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
bis (2-Cloroethyl)Ether	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2-Chlorophenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
1,3-Dichlorobenzene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
1,4-Dichlorobenzene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Benzyl Alcohol	390	U	UJ	360	U	UJ	380	U	UJ	380	U	UJ	350	U	UJ	340	U	UJ	360	U	UJ
1,2-Dichlorobenzene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2-Methylphenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
bis (2-Chloroisopropyl)Ether	390	U	UJ	360	U	UJ	380	U	UJ	380	U	---	350	U	---	340	U	---	360	U	---
4-Methylphenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
N-Nitroso-Di-n-Propylamine	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Hexachloroethane	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Nitrobenzene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Isophorone	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2-Nitrophenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2,4-Dimethylphenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Benzoic Acid	1900	U	---	1700	U	---	1900	U	---	200	J	---	1700	U	---	1700	U	---	1700	U	---
bis (2-Chloroethoxy)Methane	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2,4-Dichlorophenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
1,2,4-Trichlorobenzene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Naphthalene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
4-Chloroaniline	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Hexachlorobutadiene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**	SA2-6D-SS	Q*	DV**
4-Chloro-3-Methylphenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2-Methylnaphthalene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Hexachlorocyclopentadiene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2,4,6-Trichlorophenol	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2,4,5-Trichlorophenol	1900	U	---	1700	U	---	1900	U	---	1900	U	---	1700	U	---	1700	U	---	1700	U	---
2-Chloronaphthalene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2-Nitroaniline	1900	U	---	1700	U	---	1900	U	---	1900	U	---	1700	U	---	1700	U	---	1700	U	---
Dimethyl Phthalate	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Acenaphthylene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2,6-Dinitrotoluene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
3-Nitroaniline	1900	U	---	1700	U	---	1900	U	---	1900	U	---	1700	U	---	1700	U	---	1700	U	---
Acenaphthene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2,4-Dinitrophenol	1900	U	---	1700	U	---	1900	U	---	1900	U	UJ	1700	U	UJ	1700	U	UJ	1700	U	UJ
4-Nitrophenol	1900	U	---	1700	U	---	1900	U	---	1900	U	---	1700	U	---	1700	U	---	1700	U	---
Dibenzofuran	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
2,4-Dinitrotoluene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Diethylphthalate	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
4-Chlorophenyl-phenylether	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Fluorene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
4-Nitroaniline	1900	U	---	1700	U	---	1900	U	---	1900	U	---	1700	U	---	1700	U	---	1700	U	---
4,6-Dinitro-2-Methylphenol	1900	U	---	1700	U	---	1900	U	---	1900	U	---	1700	U	---	1700	U	---	1700	U	---
N-Nitrosodiphenylamine (1)	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
4-Bromophenyl-phenylether	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA2-5C-SS	Q*	DV**	SA2-5D-SS	Q*	DV**	SA2-5E-SS	Q*	DV**	SA2-6A-SS	Q*	DV**	SA2-6B-SS	Q*	DV**	SA2-6C-SS	Q*	DV**	SA2-6D-SS	Q*	DV**
Hexachlorobenzene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	--
Pentachlorophenol	1900	U	---	1700	U	---	1900	U	---	1900	U	---	1700	U	---	1700	U	---	1700	U	---
Phenanthrene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Anthracene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Di-n-Butylphthalate	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Fluoranthrene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Pyrene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Butylbenzylphthalate	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	--
3,3'-Dichlorbenzidine	780	U	---	720	U	---	760	U	---	770	U	---	710	U	---	690	U	---	710	U	---
Benzo(a)Anthracene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Chrysene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Bis (2-Ethylhexyl)Phthalate	64	J	---	360	U	---	50	J	---	380	U	---	350	U	---	340	U	---	360	U	---
Di-n-Octyl Phthalate	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Benzo(b)Fluoranthene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Benzo(k)Fluoranthene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Benzo(a)Pyrene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Indeno(1,2,3-cd)Pyrene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	---
Dibenz(a,h)Anthracene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	--
Benzo(g,h,i)Perylene	390	U	---	360	U	---	380	U	---	380	U	---	350	U	---	340	U	---	360	U	- -

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**	SA3-8A-SS	Q*	DV**
Phenol	360	U	---	380	U	---	400	U	---	360	U	---	48	J	---	350	U	---	400	U	---
bis (2-Cloroethyl)Ether	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2-Chlorophenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
1,3-Dichlorobenzene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
1,4-Dichlorobenzene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Benzyl Alcohol	360	U	---	380	U	UJ	400	U	UJ	360	U	UJ	360	U	UJ	350	U	---	400	U	UJ
1,2-Dichlorobenzene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2-Methylphenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
bis (2-Chloroisopropyl)Ether	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
4-Methylphenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
N-Nitroso-Di-n-Propylamine	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Hexachloroethane	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Nitrobenzene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Isophorone	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2-Nitrophenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2,4-Dimethylphenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Benzoic Acid	450	J	---	240	J	---	260	J	---	230	J	---	200	J	---	110	J	---	1900	U	---
bis (2-Chloroethoxy)Methane	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2,4-Dichlorophenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
1,2,4-Trichlorobenzene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Naphthalene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
4-Chloroaniline	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Hexachlorobutadiene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**	SA3-8A-SS	Q*	DV**
4-Chloro-3-Methylphenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2-Methylnaphthalene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Hexachlorocyclopentadiene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2,4,6-Trichlorophenol	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2,4,5-Trichlorophenol	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
2-Chloronaphthalene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2-Nitroaniline	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
Dimethyl Phthalate	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Acenaphthylene	360	U	---	380	U	---	400	U	---	360	U	---	110	J	---	350	U	---	400	U	---
2,6-Dinitrotoluene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
3-Nitroaniline	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
Acenaphthene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2,4-Dinitrophenol	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
4-Nitrophenol	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
Dibenzofuran	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
2,4-Dinitrotoluene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Diethylphthalate	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
4-Chlorophenyl-phenylether	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Fluorene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
4-Nitroaniline	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
4,6-Dinitro-2-Methylphenol	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
N-Nitrosodiphenylamine (1)	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
4-Bromophenyl-phenylether	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA2-6E-SS	Q*	DV**	SA3-7A-SS	Q*	DV**	SA3-7B-SS	Q*	DV**	SA3-7C-SS	Q*	DV**	SA3-7D-SS	Q*	DV**	SA3-7E-SS	Q*	DV**	SA3-8A-SS	Q*	DV**
Hexachlorobenzene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Pentachlorophenol	1800	U	---	1900	U	---	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1900	U	---
Phenanthrene	360	U	---	380	U	---	400	U	---	360	U	---	52	J	---	350	U	---	400	U	---
Anthracene	360	U	---	380	U	---	400	U	---	360	U	---	160	J	---	350	U	---	400	U	---
Di-n-Butylphthalate	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Fluoranthrene	360	U	---	380	U	---	400	U	---	360	U	---	2300	D	---	350	U	---	400	U	---
Pyrene	360	U	---	380	U	---	400	U	---	360	U	---	3000	D	DJ	350	U	---	400	U	---
Butylbenzylphthalate	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
3,3'-Dichlorbenzidine	720	U	---	760	U	---	790	U	---	720	U	---	720	U	---	690	U	---	800	U	---
Benzo(a)Anthracene	360	U	---	380	U	---	400	U	---	360	U	---	1200		J	350	U	---	400	U	---
Chrysene	360	U	---	380	U	---	400	U	---	360	U	---	1800		J	350	U	---	400	U	---
Bis (2-Ethylhexyl)Phthalate	130	J	---	380	U	---	400	U	---	360	U	---	360	U	---	56	J	---	400	U	---
Di-n-Octyl Phthalate	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Benzo(b)Fluoranthene	360	U	---	380	U	---	400	U	---	360	U	---	1600		---	350	U	---	400	U	---
Benzo(k)Fluoranthene	360	U	---	380	U	---	400	U	---	360	U	---	350	J	---	350	U	---	400	U	---
Benzo(a)Pyrene	360	U	---	380	U	---	400	U	---	360	U	---	520		---	350	U	---	400	U	---
Indeno(1,2,3-cd)Pyrene	360	U	---	380	U	---	400	U	---	360	U	---	110	J	---	350	U	---	400	U	---
Dibenz(a,h)Anthracene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---
Benzo(g,h,i)Perylene	360	U	---	380	U	---	400	U	---	360	U	---	360	U	---	350	U	---	400	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**	SA3-9C-SS	Q*	DV**
Phenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
bis (2-Cloroethyl)Ether	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2-Chlorophenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
1,3-Dichlorobenzene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
1,4-Dichlorobenzene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Benzyl Alcohol	350	U	UJ	340	U	UJ	350	U	UJ	390	U	UJ	410	U	UJ	390	U	UJ	360	U	UJ
1,2-Dichlorobenzene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2-Methylphenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
bis (2-Chloroisopropyl)Ether	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
4-Methylphenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
N-Nitroso-Di-n-Propylamine	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Hexachloroethane	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Nitrobenzene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Isophorone	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2-Nitrophenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2,4-Dimethylphenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Benzoic Acid	150	J	---	180	J	---	170	J	---	300	J	---	2000	U	---	260	J	---	1700	U	---
bis (2-Chloroethoxy)Methane	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2,4-Dichlorophenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
1,2,4-Trichlorobenzene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Naphthalene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
4-Chloroaniline	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Hexachlorobutadiene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**	SA3-9C-SS	Q*	DV**
4-Chloro-3-Methylphenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2-Methylnaphthalene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Hexachlorocyclopentadiene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2,4,6-Trichlorophenol	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2,4,5-Trichlorophenol	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
2-Chloronaphthalene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2-Nitroaniline	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
Dimethyl Phthalate	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Acenaphthylene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2,6-Dinitrotoluene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
3-Nitroaniline	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
Acenaphthene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2,4-Dinitrophenol	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
4-Nitrophenol	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
Dibenzofuran	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
2,4-Dinitrotoluene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Diethylphthalate	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
4-Chlorophenyl-phenylether	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Fluorene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
4-Nitroaniline	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
4,6-Dinitro-2-Methylphenol	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
N-Nitrosodiphenylamine (1)	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
4-Bromophenyl-phenylether	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA3-8B-SS	Q*	DV**	SA3-8C-SS	Q*	DV**	SA3-8D-SS	Q*	DV**	SA3-8E-SS	Q*	DV**	SA3-9A-SS	Q*	DV**	SA3-9B-SS	Q*	DV**	SA3-9C-SS	Q*	DV**
Hexachlorobenzene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Pentachlorophenol	1700	U	---	1600	U	---	1700	U	---	1900	U	---	2000	U	---	1900	U	---	1700	U	---
Phenanthrene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Anthracene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Di-n-Butylphthalate	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Fluoranthrene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Pyrene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Butylbenzylphthalate	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
3,3'-Dichlorbenzidine	700	U	---	680	U	---	690	U	---	770	U	---	820	U	---	780	U	---	720	U	---
Benzo(a)Anthracene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Chrysene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Bis (2-Ethylhexyl)Phthalate	140	J	---	340	U	---	200	J	---	390	U	---	410	U	---	390	U	---	48	J	---
Di-n-Octyl Phthalate	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Benzo(b)Fluoranthene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Benzo(k)Fluoranthene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Benzo(a)Pyrene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Indeno(1,2,3-cd)Pyrene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Dibenz(a,h)Anthracene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---
Benzo(g,h,i)Perylene	350	U	---	340	U	---	350	U	---	390	U	---	410	U	---	390	U	---	360	U	---

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**	SA3-10E-SS	Q*	DV**
Phenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
bis (2-Cloroethyl)Ether	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2-Chlorophenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
1,3-Dichlorobenzene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
1,4-Dichlorobenzene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Benzyl Alcohol	410	U	UJ	410	U	UJ	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
1,2-Dichlorobenzene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2-Methylphenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
bis (2-Chloroisopropyl)Ether	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
4-Methylphenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
N-Nitroso-Di-n-Propylamine	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Hexachloroethane	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Nitrobenzene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Isophorone	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2-Nitrophenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2,4-Dimethylphenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Benzoic Acid	120	J	---	140	J	---	870	J	---	210	J	---	850	J	---	340	J	---	460	J	---
bis (2-Chloroethoxy)Methane	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2,4-Dichlorophenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
1,2,4-Trichlorobenzene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Naphthalene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
4-Chloroaniline	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Hexachlorobutadiene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**	SA3-10E-SS	Q*	DV**
4-Chloro-3-Methylphenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2-Methylnaphthalene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Hexachlorocyclopentadiene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2,4,6-Trichlorophenol	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2,4,5-Trichlorophenol	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
2-Chloronaphthalene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2-Nitroaniline	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
Dimethyl Phthalate	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Acenaphthylene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2,6-Dinitrotoluene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
3-Nitroaniline	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
Acenaphthene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2,4-Dinitrophenol	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
4-Nitrophenol	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
Dibenzofuran	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
2,4-Dinitrotoluene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Diethylphthalate	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
4-Chlorophenyl-phenylether	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Fluorene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
4-Nitroaniline	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
4,6-Dinitro-2-Methylphenol	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
N-Nitrosodiphenylamine (1)	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
4-Bromophenyl-phenylether	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA3-9D-SS	Q*	DV**	SA3-9E-SS	Q*	DV**	SA3-10A-SS	Q*	DV**	SA3-10B-SS	Q*	DV**	SA3-10C-SS	Q*	DV**	SA3-10D-SS	Q*	DV**	SA3-10E-SS	Q*	DV**
Hexachlorobenzene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Pentachlorophenol	2000	U	---	2000	U	---	1800	U	---	1900	U	---	1800	U	---	1800	U	---	2000	U	---
Phenanthrene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Anthracene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Di-n-Butylphthalate	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Fluoranthrene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Pyrene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Butylbenzylphthalate	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
3,3'-Dichlorbenzidine	820	U	---	820	U	---	750	U	---	800	U	---	760	U	---	750	U	---	810	U	---
Benzo(a)Anthracene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Chrysene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Bis (2-Ethylhexyl)Phthalate	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Di-n-Octyl Phthalate	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Benzo(b)Fluoranthene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Benzo(k)Fluoranthene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Benzo(a)Pyrene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Indeno(1,2,3-cd)Pyrene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Dibenz(a,h)Anthracene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---
Benzo(g,h,i)Perylene	410	U	---	410	U	---	370	U	---	400	U	---	380	U	---	370	U	---	400	U	---

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA4-11A-SS	Q*	DV**	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**	SA4-12B-SS	Q*	DV**
Phenol	340	U	---	130	J	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
bis (2-Cloroethyl)Ether	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2-Chlorophenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
1,3-Dichlorobenzene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
1,4-Dichlorobenzene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Benzyl Alcohol	340	U	J	340	U	J	340	U	J	350	U	J	330	U	J	400	U	---	370	U	---
1,2-Dichlorobenzene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2-Methylphenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
bis (2-Chloroisopropyl)Ether	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
4-Methylphenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
N-Nitroso-Di-n-Propylamine	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Hexachloroethane	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Nitrobenzene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Isophorone	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2-Nitrophenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2,4-Dimethylphenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Benzoic Acid	1600	U	---	1600	U	---	75	J	---	820	J	---	1600	U	---	89	J	---	120	J	---
bis (2-Chloroethoxy)Methane	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2,4-Dichlorophenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
1,2,4-Trichlorobenzene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Naphthalene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
4-Chloroaniline	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Hexachlorobutadiene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA4-11A-SS	Q*	DV**	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**	SA4-12B-SS	Q*	DV**
4-Chloro-3-Methylphenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2-Methylnaphthalene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Hexachlorocyclopentadiene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2,4,6-Trichlorophenol	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2,4,5-Trichlorophenol	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
2-Chloronaphthalene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2-Nitroaniline	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
Dimethyl Phthalate	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Acenaphthylene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2,6-Dinitrotoluene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
3-Nitroaniline	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
Acenaphthene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2,4-Dinitrophenol	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
4-Nitrophenol	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
Dibenzofuran	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
2,4-Dinitrotoluene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Diethylphthalate	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
4-Chlorophenyl-phenylether	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Fluorene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
4-Nitroaniline	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
4,6-Dinitro-2-Methylphenol	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
N-Nitrosodiphenylamine (1)	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
4-Bromophenyl-phenylether	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA4-11A-SS	Q*	DV**	SA4-11B-SS	Q*	DV**	SA4-11C-SS	Q*	DV**	SA4-11D-SS	Q*	DV**	SA4-11E-SS	Q*	DV**	SA4-12A-SS	Q*	DV**	SA4-12B-SS	Q*	DV**
Hexachlorobenzene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Pentachlorophenol	1600	U	---	1600	U	---	1600	U	---	1700	U	---	1600	U	---	1900	U	---	1800	U	---
Phenanthrene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Anthracene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Di-n-Butylphthalate	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Fluoranthrene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Pyrene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Butylbenzylphthalate	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
3,3'-Dichlorbenzidine	670	U	---	670	U	---	680	U	---	700	U	---	670	U	---	800	U	---	730	U	---
Benzo(a)Anthracene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Chrysene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Bis (2-Ethylhexyl)Phthalate	62	J	---	55	J	---	54	J	---	51	J	---	81	J	---	400	U	---	100	J	---
Di-n-Octyl Phthalate	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Benzo(b)Fluoranthene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Benzo(k)Fluoranthene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Benzo(a)Pyrene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Indeno(1,2,3-cd)Pyrene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Dibenz(a,h)Anthracene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	---	370	U	---
Benzo(g,h,i)Perylene	340	U	---	340	U	---	340	U	---	350	U	---	330	U	---	400	U	UJ	370	U	UJ

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**	SA4-13D-SS	Q*	DV**
Phenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
bis (2-Cloroethyl)Ether	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2-Chlorophenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
1,3-Dichlorobenzene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
1,4-Dichlorobenzene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Benzyl Alcohol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
1,2-Dichlorobenzene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2-Methylphenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
bis (2-Chloroisopropyl)Ether	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
4-Methylphenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
N-Nitroso-Di-n-Propylamine	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Hexachloroethane	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Nitrobenzene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Isophorone	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2-Nitrophenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2,4-Dimethylphenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Benzoic Acid	160	J	---	210	J	---	150	J	---	1700	U	---	1700	U	---	1700	U	NP	150	J	---
bis (2-Chloroethoxy)Methane	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2,4-Dichlorophenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
1,2,4-Trichlorobenzene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Naphthalene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
4-Chloroaniline	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Hexachlorobutadiene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**	SA4-13D-SS	Q*	DV**
4-Chloro-3-Methylphenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2-Methylnaphthalene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Hexachlorocyclopentadiene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2,4,6-Trichlorophenol	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2,4,5-Trichlorophenol	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
2-Chloronaphthalene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2-Nitroaniline	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
Dimethyl Phthalate	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Acenaphthylene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2,6-Dinitrotoluene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
3-Nitroaniline	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
Acenaphthene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2,4-Dinitrophenol	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
4-Nitrophenol	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
Dibenzofuran	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
2,4-Dinitrotoluene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Diethylphthalate	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
4-Chlorophenyl-phenylether	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Fluorene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
4-Nitroaniline	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
4,6-Dinitro-2-Methylphenol	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
N-Nitrosodiphenylamine (1)	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
4-Bromophenyl-phenylether	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA4-12C-SS	Q*	DV**	SA4-12D-SS	Q*	DV**	SA4-12E-SS	Q*	DV**	SA4-13A-SS	Q*	DV**	SA4-13B-SS	Q*	DV**	SA4-13C-SS	Q*	DV**	SA4-13D-SS	Q*	DV**
Hexachlorobenzene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Pentachlorophenol	1800	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	NP	1800	U	---
Phenanthrene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Anthracene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Di-n-Butylphthalate	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Fluoranthrene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Pyrene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Butylbenzylphthalate	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
3,3'-Dichlorbenzidine	760	U	---	760	U	---	690	U	---	690	U	---	710	U	---	690	U	NP	730	U	---
Benzo(a)Anthracene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Chrysene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Bis (2-Ethylhexyl)Phthalate	79	J	---	380	U	---	350	U	---	36	J	---	350	U	---	77	J	NP	54	J	---
Di-n-Octyl Phthalate	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Benzo(b)Fluoranthene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Benzo(k)Fluoranthene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Benzo(a)Pyrene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Indeno(1,2,3-cd)Pyrene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Dibenz(a,h)Anthracene	380	U	---	380	U	---	350	U	---	340	U	---	350	U	---	340	U	NP	370	U	---
Benzo(g,h,i)Perylene	380	U	UJ	380	U	UJ	350	U	UJ	340	U	UJ	350	U	UJ	340	U	NP	370	U	UJ

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**	SA4-15A-SS	Q*	DV**
Phenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
bis (2-Cloroethyl)Ether	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2-Chlorophenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
1,3-Dichlorobenzene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
1,4-Dichlorobenzene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Benzyl Alcohol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	UJ	350	U	UJ	340	U	---
1,2-Dichlorobenzene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2-Methylphenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
bis (2-Chloroisopropyl)Ether	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
4-Methylphenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
N-Nitroso-Di-n-Propylamine	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Hexachloroethane	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Nitrobenzene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Isophorone	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2-Nitrophenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2,4-Dimethylphenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Benzoic Acid	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
bis (2-Chloroethoxy)Methane	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2,4-Dichlorophenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
1,2,4-Trichlorobenzene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Naphthalene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
4-Chloroaniline	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Hexachlorobutadiene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**	SA4-15A-SS	Q*	DV**
4-Chloro-3-Methylphenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2-Methylnaphthalene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Hexachlorocyclopentadiene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2,4,6-Trichlorophenol	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2,4,5-Trichlorophenol	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
2-Chloronaphthalene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2-Nitroaniline	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
Dimethyl Phthalate	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Acenaphthylene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2,6-Dinitrotoluene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
3-Nitroaniline	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
Acenaphthene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2,4-Dinitrophenol	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	UJ	1700	U	UJ	1700	U	---
4-Nitrophenol	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
Dibenzofuran	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
2,4-Dinitrotoluene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Diethylphthalate	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
4-Chlorophenyl-phenylether	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Fluorene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
4-Nitroaniline	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
4,6-Dinitro-2-Methylphenol	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
N-Nitrosodiphenylamine (1)	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
4-Bromophenyl-phenylether	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA4-13E-SS	Q*	DV**	SA4-14A-SS	Q*	DV**	SA4-14B-SS	Q*	DV**	SA4-14C-SS	Q*	DV**	SA4-14D-SS	Q*	DV**	SA4-14E-SS	Q*	DV**	SA4-15A-SS	Q*	DV**
Hexachlorobenzene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Pentachlorophenol	1900	U	---	1700	U	---	1800	U	---	1700	U	---	1700	U	---	1700	U	---	1700	U	---
Phenanthrene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Anthracene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Di-n-Butylphthalate	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Fluoranthrene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Pyrene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Butylbenzylphthalate	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
3,3'-Dichlorbenzidine	780	U	---	690	U	---	730	U	---	720	U	---	690	U	---	700	U	---	690	U	---
Benzo(a)Anthracene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Chrysene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Bis (2-Ethylhexyl)Phthalate	43	J	---	48	J	---	68	J	---	91	J	---	57	J	---	54	J	---	340	U	---
Di-n-Octyl Phthalate	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Benzo(b)Fluoranthene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Benzo(k)Fluoranthene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Benzo(a)Pyrene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Indeno(1,2,3-cd)Pyrene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Dibenz(a,h)Anthracene	390	U	---	350	U	---	370	U	---	360	U	---	340	U	---	350	U	---	340	U	---
Benzo(g,h,i)Perylene	390	U	UJ	350	U	UJ	370	U	UJ	360	U	UJ	340	U	---	350	U	---	340	U	UJ

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**	SA5-16B-SS	Q*	DV**	SA5-16C-SS	Q*	DV**
Phenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
bis (2-Cloroethyl)Ether	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2-Chlorophenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
1,3-Dichlorobenzene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
1,4-Dichlorobenzene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Benzyl Alcohol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	UJ	360	U	---	360	U	---
1,2-Dichlorobenzene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2-Methylphenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
bis (2-Chloroisopropyl)Ether	390	U	---	370	U	---	380	U	---	380	U	---	370	U	UJ	360	U	---	360	U	---
4-Methylphenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
N-Nitroso-Di-n-Propylamine	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Hexachloroethane	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Nitrobenzene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Isophorone	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2-Nitrophenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2,4-Dimethylphenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Benzoic Acid	140	J	---	1800	U	---	1900	U	---	220	J	---	1800	U	---	1700	U	---	1700	U	---
bis (2-Chloroethoxy)Methane	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2,4-Dichlorophenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
1,2,4-Trichlorobenzene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Naphthalene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
4-Chloroaniline	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Hexachlorobutadiene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**	SA5-16B-SS	Q*	DV**	SA5-16C-SS	Q*	DV**
4-Chloro-3-Methylphenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2-Methylnaphthalene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Hexachlorocyclopentadiene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2,4,6-Trichlorophenol	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2,4,5-Trichlorophenol	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
2-Chloronaphthalene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2-Nitroaniline	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
Dimethyl Phthalate	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Acenaphthylene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2,6-Dinitrotoluene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
3-Nitroaniline	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
Acenaphthene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2,4-Dinitrophenol	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
4-Nitrophenol	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
Dibenzofuran	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
2,4-Dinitrotoluene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Diethylphthalate	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
4-Chlorophenyl-phenylether	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Fluorene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
4-Nitroaniline	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
4,6-Dinitro-2-Methylphenol	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
N-Nitrosodiphenylamine (1)	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
4-Bromophenyl-phenylether	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA4-15B-SS	Q*	DV**	SA4-15C-SS	Q*	DV**	SA4-15D-SS	Q*	DV**	SA4-15E-SS	Q*	DV**	SA5-16A-SS	Q*	DV**	SA5-16B-SS	Q*	DV**	SA5-16C-SS	Q*	DV**
Hexachlorobenzene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Pentachlorophenol	1900	U	---	1800	U	---	1900	U	---	1900	U	---	1800	U	---	1700	U	---	1700	U	---
Phenanthrene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Anthracene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Di-n-Butylphthalate	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Fluoranthrene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Pyrene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Butylbenzylphthalate	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
3,3'-Dichlorbenzidine	790	U	---	730	U	---	770	U	---	760	U	---	730	U	---	710	U	---	710	U	---
Benzo(a)Anthracene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Chrysene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Bis (2-Ethylhexyl)Phthalate	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	71	J	---	90	J	---
Di-n-Octyl Phthalate	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Benzo(b)Fluoranthene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Benzo(k)Fluoranthene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Benzo(a)Pyrene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Indeno(1,2,3-cd)Pyrene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Dibenz(a,h)Anthracene	390	U	---	370	U	---	380	U	---	380	U	---	370	U	---	360	U	---	360	U	---
Benzo(g,h,i)Perylene	390	U	UJ	370	U	UJ	380	U	UJ	380	U	UJ	370	U	---	360	U	UJ	360	U	UJ

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**	SA5-17D-SS	Q*	DV**	SA5-17E-SS	Q*	DV**
Phenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
bis (2-Cloroethyl)Ether	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2-Chlorophenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
1,3-Dichlorobenzene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
1,4-Dichlorobenzene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Benzyl Alcohol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
1,2-Dichlorobenzene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2-Methylphenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
bis (2-Chloroisopropyl)Ether	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
4-Methylphenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
N-Nitroso-Di-n-Propylamine	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Hexachloroethane	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Nitrobenzene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Isophorone	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2-Nitrophenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2,4-Dimethylphenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Benzoic Acid	69	J	---	85	J	---	180	J	---	110	J	---	1700	U	---	75	J	---	1800	U	---
bis (2-Chloroethoxy)Methane	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2,4-Dichlorophenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
1,2,4-Trichlorobenzene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Naphthalene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
4-Chloroaniline	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Hexachlorobutadiene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**	SA5-17D-SS	Q*	DV**	SA5-17E-SS	Q*	DV**
4-Chloro-3-Methylphenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2-Methylnaphthalene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Hexachlorocyclopentadiene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2,4,6-Trichlorophenol	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2,4,5-Trichlorophenol	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
2-Chloronaphthalene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2-Nitroaniline	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
Dimethyl Phthalate	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Acenaphthylene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2,6-Dinitrotoluene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
3-Nitroaniline	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
Acenaphthene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2,4-Dinitrophenol	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
4-Nitrophenol	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
Dibenzofuran	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
2,4-Dinitrotoluene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Diethylphthalate	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
4-Chlorophenyl-phenylether	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Fluorene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
4-Nitroaniline	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
4,6-Dinitro-2-Methylphenol	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
N-Nitrosodiphenylamine (1)	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
4-Bromophenyl-phenylether	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA5-16D-SS	Q*	DV**	SA5-16E-SS	Q*	DV**	SA5-17A-SS	Q*	DV**	SA5-17B-SS	Q*	DV**	SA5-17C-SS	Q*	DV**	SA5-17D-SS	Q*	DV**	SA5-17E-SS	Q*	DV**
Hexachlorobenzene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Pentachlorophenol	1700	U	---	2100	U	---	1800	U	---	1800	U	---	1700	U	---	1900	U	---	1800	U	---
Phenanthrene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Anthracene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Di-n-Butylphthalate	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Fluoranthrene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Pyrene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Butylbenzylphthalate	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
3,3'-Dichlorbenzidine	700	U	---	860	U	---	750	U	---	750	U	---	720	U	---	770	U	---	750	U	---
Benzo(a)Anthracene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Chrysene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	---
Bis (2-Ethylhexyl)Phthalate	48	J	---	67	J	---	76	J	---	63	J	---	150	J	---	61	J	---	120	J	---
Di-n-Octyl Phthalate	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	UJ
Benzo(b)Fluoranthene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	UJ
Benzo(k)Fluoranthene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	UJ
Benzo(a)Pyrene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	UJ
Indeno(1,2,3-cd)Pyrene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	UJ
Dibenz(a,h)Anthracene	350	U	---	430	U	---	370	U	---	370	U	---	360	U	---	380	U	---	370	U	UJ
Benzo(g,h,i)Perylene	350	U	UJ	430	U	UJ	370	U	UJ	370	U	UJ	360	U	UJ	380	U	UJ	370	U	UJ

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles																							
	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**	REFC-19A-SS	Q*	DV**	REFC-19B-SS	Q*	DV**			
Phenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
bis (2-Cloroethyl)Ether	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
2-Chlorophenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
1,3-Dichlorobenzene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
1,4-Dichlorobenzene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
Benzyl Alcohol	370	U	UJ	370	U	UJ	370	U	UJ	360	U	UJ	370	U	UJ	350	U	---	380	U	---			
1,2-Dichlorobenzene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
2-Methylphenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
bis (2-Chloroisopropyl)Ether	370	U	UJ	370	U	UJ	370	U	UJ	360	U	UJ	370	U	UJ	350	U	---	380	U	---			
4-Methylphenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
N-Nitroso-Di-n-Propylamine	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
Hexachloroethane	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
Nitrobenzene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
Isophorone	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
2-Nitrophenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
2,4-Dimethylphenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
Benzoic Acid	1800	U	---	1800	U	---	130	J	---	1700	U	---	1800	U	---	230	J	---	310	J	---			
bis (2-Chloroethoxy)Methane	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
2,4-Dichlorophenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
1,2,4-Trichlorobenzene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
Naphthalene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
4-Chloroaniline	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			
Hexachlorobutadiene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---			

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**	REFC-19A-SS	Q*	DV**	REFC-19B-SS	Q*	DV**
4-Chloro-3-Methylphenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
2-Methylnaphthalene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Hexachlorocyclopentadiene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
2,4,6-Trichlorophenol	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
2,4,5-Trichlorophenol	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
2-Chloronaphthalene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
2-Nitroaniline	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
Dimethyl Phthalate	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Acenaphthylene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
2,6-Dinitrotoluene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
3-Nitroaniline	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
Acenaphthene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
2,4-Dinitrophenol	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
4-Nitrophenol	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
Dibenzofuran	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
2,4-Dinitrotoluene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Diethylphthalate	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
4-Chlorophenyl-phenylether	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Fluorene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
4-Nitroaniline	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
4,6-Dinitro-2-Methylphenol	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
N-Nitrosodiphenylamine (1)	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
4-Bromophenyl-phenylether	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---

* Data Qualifier. ** Data Validation.

Table A-5.4 Semi-Volatiles

Compound (ug/Kg)	SA6-18A-SS	Q*	DV**	SA6-18B-SS	Q*	DV**	SA6-18C-SS	Q*	DV**	SA6-18D-SS	Q*	DV**	SA6-18E-SS	Q*	DV**	REFC-19A-SS	Q*	DV**	REFC-19B-SS	Q*	DV**
Hexachlorobenzene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Pentachlorophenol	1800	U	---	1800	U	---	1800	U	---	1700	U	---	1800	U	---	1700	U	---	1800	U	---
Phenanthrene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Anthracene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Di-n-Butylphthalate	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Fluoranthrene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Pyrene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Butylbenzylphthalate	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
3,3'-Dichlorbenzidine	750	U	---	740	U	---	730	U	---	720	U	---	730	U	---	710	U	---	760	U	---
Benzo(a)Anthracene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Chrysene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Bis (2-Ethylhexyl)Phthalate	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Di-n-Octyl Phthalate	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Benzo(b)Fluoranthene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Benzo(k)Fluoranthene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Benzo(a)Pyrene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Indeno(1,2,3-cd)Pyrene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Dibenz(a,h)Anthracene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	---	380	U	---
Benzo(g,h,i)Perylene	370	U	---	370	U	---	370	U	---	360	U	---	370	U	---	350	U	UJ	380	U	UJ

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles								
	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**
Phenol	450	U	---	340	U	---	350	U	---
bis (2-Cloroethyl)Ether	450	U	---	340	U	---	350	U	---
2-Chlorophenol	450	U	---	340	U	---	350	U	---
1,3-Dichlorobenzene	450	U	---	340	U	---	350	U	---
1,4-Dichlorobenzene	450	U	---	340	U	---	350	U	---
Benzyl Alcohol	450	U	---	340	U	---	350	U	---
1,2-Dichlorobenzene	450	U	---	340	U	---	350	U	---
2-Methylphenol	450	U	---	340	U	---	350	U	---
bis (2-Chloroisopropyl)Ether	450	U	---	340	U	---	350	U	---
4-Methylphenol	450	U	---	340	U	---	350	U	---
N-Nitroso-Di-n-Propylamine	450	U	---	340	U	---	350	U	---
Hexachloroethane	450	U	---	340	U	---	350	U	---
Nitrobenzene	450	U	---	340	U	---	350	U	---
Isophorone	450	U	---	340	U	---	350	U	---
2-Nitrophenol	450	U	---	340	U	---	350	U	---
2,4-Dimethylphenol	450	U	---	340	U	---	350	U	---
Benzoic Acid	570	J	---	99	J	---	1700	U	---
bis (2-Chloroethoxy)Methane	450	U	---	340	U	---	350	U	---
2,4-Dichlorophenol	450	U	---	340	U	---	350	U	---
1,2,4-Trichlorobenzene	450	U	---	340	U	—	350	U	—
Naphthalene	450	U	---	340	U	—	350	U	---
4-Chloroaniline	450	U	—	340	U	—	350	U	—
Hexachlorobutadiene	450	U	---	340	U	—	350	U	—

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles								
	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**
4-Chloro-3-Methylphenol	450	U	---	340	U	---	350	U	---
2-Methylnaphthalene	450	U	---	340	U	---	350	U	---
Hexachlorocyclopentadiene	450	U	---	340	U	---	350	U	---
2,4,6-Trichlorophenol	450	U	---	340	U	---	350	U	---
2,4,5-Trichlorophenol	2200	U	---	1600	U	---	1700	U	---
2-Chloronaphthalene	450	U	---	340	U	---	350	U	---
2-Nitroaniline	2200	U	---	1600	U	---	1700	U	---
Dimethyl Phthalate	450	U	---	340	U	---	350	U	---
Acenaphthylene	450	U	---	340	U	---	350	U	---
2,6-Dinitrotoluene	450	U	---	340	U	---	350	U	---
3-Nitroaniline	2200	U	---	1600	U	---	1700	U	---
Acenaphthene	450	U	---	340	U	---	350	U	---
2,4-Dinitrophenol	2200	U	---	1600	U	---	1700	U	---
4-Nitrophenol	2200	U	---	1600	U	---	1700	U	---
Dibenzofuran	450	U	---	340	U	---	350	U	---
2,4-Dinitrotoluene	450	U	---	340	U	---	350	U	---
Diethylphthalate	450	U	---	340	U	---	350	U	---
4-Chlorophenyl-phenylether	450	U	---	340	U	---	350	U	---
Fluorene	450	U	---	340	U	---	350	U	---
4-Nitroaniline	2200	U	---	1600	U	---	1700	U	---
4,6-Dinitro-2-Methylphenol	2200	U	---	1600	U	---	1700	U	---
N-Nitrosodiphenylamine (1)	450	U	---	340	U	---	350	U	---
4-Bromophenyl-phenylether	450	U	---	340	U	---	350	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-5.4 Semi-Volatiles								
	REFC-19C-SS	Q*	DV**	REFC-19D-SS	Q*	DV**	REFC-19E-SS	Q*	DV**
Hexachlorobenzene	450	U	---	340	U	---	350	U	---
Pentachlorophenol	2200	U	---	1600	U	---	1700	U	---
Phenanthrene	450	U	---	340	U	---	350	U	---
Anthracene	450	U	---	340	U	---	350	U	---
Di-n-Butylphthalate	450	U	---	340	U	---	350	U	---
Fluoranthrene	450	U	---	340	U	---	350	U	---
Pyrene	450	U	---	340	U	---	350	U	---
Butylbenzylphthalate	450	U	---	340	U	---	350	U	---
3,3'-Dichlorbenzidine	900	U	---	680	U	---	700	U	---
Benzo(a)Anthracene	450	U	---	340	U	---	350	U	---
Chrysene	450	U	---	340	U	---	350	U	---
Bis (2-Ethylhexyl)Phthalate	450	U	---	340	U	---	350	U	---
Di-n-Octyl Phthalate	450	U	---	340	U	---	350	U	---
Benzo(b)Fluoranthene	450	U	---	340	U	---	350	U	---
Benzo(k)Fluoranthene	450	U	---	340	U	---	350	U	---
Benzo(a)Pyrene	450	U	---	340	U	---	350	U	---
Indeno(1,2,3-cd)Pyrene	450	U	---	340	U	---	350	U	---
Dibenz(a,h)Anthracene	450	U	---	340	U	---	350	U	---
Benzo(g,h,i)Perylene	450	U	UJ	340	U	UJ	350	U	UJ

Parameter (pCi/g)	Table A-5.5 Gross Alpha/Beta																			
	SA1-1A-SS	DV**	SA1-1A-SS-D	DV**	SA1-1B-SS	DV**	SA1-1C-SS	DV**	SA1-1D-SS	DV**	SA1-1D-SS-D	DV**	SA1-1E-SS	DV**	SA1-2A-SS	DV**	SA1-2B-SS	DV**
Alpha	7.97	---	ND	---	ND	---	ND	---	ND	---	ND	---	7.98	---	7.56	---	ND	---
Sigma Error (+/-)	5.24	---	-	---	-	---	-	---	-	---	-	---	5.28	---	4.75	---	-	---
MDA/LLD*	7.49	---	9.95	U	8.69	U	12.40	U	8.90	U	6.46	U	7.73	---	6.59	---	7.04	U
Beta	15.70	---	16.90	---	9.74	---	9.88	---	11.40	---	10.90	---	10.10	---	8.97	---	12.10	---
Sigma Error (+/-)	3.69	---	4.81	---	3.64	---	3.61	---	3.35	---	3.05	---	3.21	---	3.24	---	3.32	---
MDA/LLD*	4.62	---	6.65	---	5.37	---	5.27	---	4.54	---	4.15	---	4.57	---	4.75	---	4.54	---

Table A-5.5 Gross Alpha/Beta																				
Parameter (pCi/g)	SA1-2C-SS	DV**	SA1-2D-SS	DV**	SA1-2E-SS	DV**	SA1-3A-SS	DV**	SA1-3B-SS	DV**	SA1-3C-SS	DV**	SA1-3D-SS	DV**	SA1-3E-SS	DV**	SA1-4A-SS	DV**
Alpha	ND	---	5.52	---	10.50	---	10.00	---	10.90	---	7.05	---	10.70	---	ND	---	11.90	---
Sigma Error (+/-)	-	---	3.86	---	4.97	---	4.37	---	4.92	---	4.93	---	5.77	---	-	---	5.34	---
MDA/LLD*	5.79	U	5.47	---	5.94	---	4.72	---	5.04	---	6.99	---	7.71	---	9.82	U	6.57	---
Beta	10.70	---	10.90	---	13.70	---	12.30	---	10.80	---	9.51	---	14.30	---	9.84	---	18.00	---
Sigma Error (+/-)	3.13	---	3.04	---	3.50	---	3.08	---	4.16	---	3.43	---	3.66	---	4.30	---	3.63	---
MDA/LLD*	4.31	---	4.12	---	4.62	---	4.02	---	6.28	---	4.96	---	4.75	---	6.56	---	4.22	---

* Minimum Detectable Activity. Lower Level of Detection. ** Data Validation.

Parameter (pCi/g)	Table A-5.5 Gross Alpha/Beta																			
	SA1-4A-SS-D	DV**	SA1-4B-SS	DV**	SA1-4C-SS	DV**	SA1-4D-SS	DV**	SA1-4E-SS	DV**	SA2-5A-SS	DV**	SA2-5A-SS-D	DV**	SA2-5B-SS	DV**	SA2-5C-SS	DV**
Alpha	8.23	---	6.85	---	ND	---	ND	---	6.70	---	13.80	---	14.70	---	18.70	---	19.90	---
Sigma Error (+/-)	5.45	---	4.30	---	-	---	-	---	4.30	---	5.72	---	5.94	---	7.10	---	7.35	---
MDA/LLD*	7.97	---	5.96	---	7.77	---	5.93	---	5.90	---	6.36	---	6.92	---	8.18	---	8.34	---
Beta	18.40	---	13.90	---	9.21	---	11.90	---	13.50	---	19.20	---	16.70	---	20.20	---	18.40	---
Sigma Error (+/-)	3.90	---	3.35	---	3.39	---	3.27	---	3.33	---	3.82	---	3.49	---	4.41	---	4.03	---
MDA/LLD*	4.72	---	4.30	---	5.01	---	4.41	---	4.45	---	4.31	---	4.08	---	5.47	---	4.92	---

* Minimum Detectable Activity. Lower Level of Detection. ** Data Validation.

Table A-5.5 Gross Alpha/Beta																								
Parameter (pCi/g)	SA2-5D-SS	DV**		SA2-5E-SS	DV**		SA2-6A-SS	DV**		SA2-6B-SS	DV**		SA2-6C-SS	DV**		SA2-6D-SS	DV**		SA2-6E-SS	DV**		SA3-7A-SS	DV**		SA3-7B-SS	DV**	
Alpha	17.80	---		17.70	---		12.60	---		6.17	---		6.16	---		8.55	---		13.80	---		7.00	---		7.54	---	
Sigma Error (+/-)	8.92	---		7.12	---		4.94	---		3.97	---		3.96	---		4.06	---		5.15	---		4.18	---		3.99	---	
MDA/LLD*	12.60	---		8.46	---		5.27	---		5.44	---		5.43	---		4.64	---		4.59	---		5.56	---		4.83	---	
Beta	20.40	---		18.00	---		17.00	---		10.50	---		11.80	---		9.98	---		8.73	---		13.00	---		11.20	---	
Sigma Error (+/-)	4.43	---		3.78	---		3.31	---		2.99	---		3.04	---		2.87	---		3.72	---		3.29	---		3.06	---	
MDA/LLD*	5.34	---		4.37	---		3.74	---		4.10	---		4.03	---		3.95	---		5.71	---		4.33	---		4.12	---	

Table A-5.5 Gross Alpha/Beta

Parameter (pCi/g)	SA3-7C-SS	DV**	SA3-7D-SS	DV**	SA3-7E-SS	DV**	SA3-8A-SS	DV**	SA3-8B-SS	DV**	SA3-8C-SS	DV**	SA3-8D-SS	DV**	SA3-8E-SS	DV**	SA3-9A-SS	DV**
Alpha	9.98	---	12.20	---	ND	---	ND	---	ND	---	ND	---	10.70	---	ND	---	37.90	J
Sigma Error (+/-)	4.49	---	5.79	---	-	---	-	---	-	---	-	---	6.03	---	-	---	12.70	---
MDA/LLD*	4.61	---	6.91	---	7.18	---	8.27	---	8.41	---	12.27	---	8.29	---	5.86	---	9.00	---
Beta	9.45	---	12.80	---	13.90	---	10.50	---	15.40	---	6.78	---	22.20	---	12.50	---	14.20	J
Sigma Error (+/-)	3.78	---	3.63	---	3.95	---	3.76	---	3.95	---	3.38	---	4.18	---	3.29	---	5.10	---
MDA/LLD*	5.74	---	4.91	---	4.43	---	5.53	---	5.20	---	5.22	---	4.49	---	4.36	---	5.96	---

* Minimum Detectable Activity, Lower Level of Detection. ** Data Validation.

Table A-5.5 Gross Alpha/Beta																				
Parameter (pCi/g)	SA3-9B-SS	DV**	SA3-9C-SS	DV**	SA3-9D-SS	DV**	SA3-9E-SS	DV**	SA3-10A-SS	DV**	SA3-10B-SS	DV**	SA3-10C-SS	DV**	SA3-10D-SS	DV**	SA3-10D-SS-D	DV**
Alpha	17.50	J	12.30	J	24.40	J	13.00	J	22.50	J	16.80	J	16.50	J	16.40	J	20.50	J
Sigma Error (+/-)	8.30	---	7.70	---	10.40	---	7.80	---	9.80	---	8.30	---	8.50	---	8.00	---	9.30	---
MDA/LLD*	8.40	---	9.89	---	1.05	---	10.00	---	9.95	---	9.18	---	9.62	---	8.15	---	9.45	---
Beta	17.80	J	13.00	J	15.50	J	14.60	J	16.00	J	11.70	J	14.40	J	9.70	J	10.10	J
Sigma Error (+/-)	5.80	---	5.30	---	5.30	---	5.30	---	5.20	---	4.50	---	5.10	---	4.39	---	4.60	---
MDA/LLD*	6.21	---	6.66	---	6.07	---	6.23	---	5.62	---	5.52	---	5.94	---	5.83	---	6.04	---

* Minimum Detectable Activity. Lower Level of Detection. ** Data Validation.

Parameter (pCi/g)	Table A-5.5 Gross Alpha/Beta																					
	SA3-10E-SS	DV**	SA4-11A-SS	DV**	SA4-11B-SS	DV**	SA4-11C-SS	DV**	SA4-11D-SS	DV**	SA4-11E-SS	DV**	SA4-12A-SS	DV**	SA4-12A-SS-D	DV**	SA4-12B-SS	DV**				
Alpha	15.80	J	6.81	---	5.66	---	7.99	---	12.30	---	9.68	---	15.70	---	20.00	---	17.90	---				
Sigma Error (+/-)	8.50	---	4.37	---	3.96	---	3.99	---	5.10	---	5.40	---	8.64	---	7.56	---	5.88	---				
MDA/LLD*	10.02	---	6.00	---	5.61	---	4.69	---	4.93	---	6.98	---	12.30	---	8.45	---	4.68	---				
Beta	11.50	J	14.20	---	11.20	---	11.80	---	16.40	---	13.60	---	26.20	---	21.50	---	31.20	---				
Sigma Error (+/-)	4.70	---	3.50	---	3.18	---	3.04	---	4.44	---	3.72	---	4.85	---	4.07	---	5.25	---				
MDA/LLD*	6.01	---	4.52	---	4.37	---	3.99	---	6.14	---	4.96	---	5.24	---	4.37	---	5.56	---				

Parameter (pCi/g)	Table A-5.5 Gross Alpha/Beta																			
	SA4-12C-SS	DV**	SA4-12D-SS	DV**	SA4-12E-SS	DV**	SA4-13A-SS	DV**	SA4-13B-SS	DV**	SA4-13C-SS	DV**	SA4-13D-SS	DV**	SA4-13E-SS	DV**	SA4-14A-SS	DV**
Alpha	11.70	---	14.90	---	ND	---	8.63	---	8.69	---	7.52	---	ND	---	ND	---	9.75	---
Sigma Error (+/-)	5.32	---	6.03	---	-	---	4.30	---	4.85	---	4.94	---	-	---	-	---	4.63	---
MDA/LLD*	6.20	---	7.02	--	8.44	U	4.71	---	6.26	---	7.07	---	8.47	U	8.12	U	5.52	---
Beta	14.50	---	14.80	---	12.80	---	13.00	---	13.40	---	11.60	---	9.26	---	14.50	---	11.60	---
Sigma Error (+/-)	3.49	---	3.36	---	3.97	---	3.92	---	3.44	---	3.14	---	3.61	---	3.78	---	3.18	---
MDA/LLD*	4.40	---	4.13	---	5.64	---	5.60	---	4.45	---	4.17	---	5.41	---	5.02	---	4.30	---

Table A-5.5 Gross Alpha/Beta																						
Parameter (pCi/g)	SA4-14A-SS-D	DV**	SA4-14B-SS	DV**	SA4-14C-SS	DV**	SA4-14D-SS	DV**	SA4-14E-SS	DV**	SA4-15A-SS	DV**	SA4-15A-SS-D	DV**	SA4-15B-SS	DV**	SA4-15C-SS	DV**
Alpha	8.03	---	10.40	---	12.10	---	ND	---	ND	---	7.54	---	16.90	---	20.50	---	11.10	---
Sigma Error (+/-)	4.01	---	4.67	---	5.48	---	-	---	-	---	4.20	---	5.52	---	6.19	---	4.64	---
MDA/LLD*	4.72	---	4.79	---	6.38	---	11.40	U	8.35	U	5.44	---	5.16	---	5.43	---	4.90	---
Beta	11.00	---	13.70	---	11.00	---	9.56	---	12.20	---	15.70	---	13.90	---	17.70	---	13.90	---
Sigma Error (+/-)	2.98	---	4.30	---	3.18	---	3.63	---	3.29	---	3.27	---	3.14	---	3.50	---	3.29	---
MDA/LLD*	4.01	---	5.97	---	4.33	---	4.86	---	4.31	---	3.86	---	3.89	---	4.04	---	4.17	---

Parameter (pCi/g)	Table A-5.5 Gross Alpha/Beta																			
	SA4-15D-SS	DV**	SA4-15E-SS	DV**	SA5-16A-SS	DV**	SA5-16B-SS	DV**	SA5-16C-SS	DV**	SA5-16D-SS	DV**	SA5-16E-SS	DV**	SA5-17A-SS	DV**	SA5-17B-SS	DV**
Alpha	12.80	---	16.80	---	18.10	---	11.00	---	12.20	---	9.11	---	14.80	---	ND	---	18.80	---
Sigma Error (+/-)	5.10	---	6.18	---	6.92	---	5.37	---	6.19	---	5.81	---	8.04	---	-	---	8.44	---
MDA/LLD*	4.79	---	6.33	---	4.71	---	6.88	---	8.18	---	8.33	---	11.60	---	8.17	U	11.50	---
Beta	23.10	---	17.10	---	23.80	---	16.10	---	16.50	---	13.30	---	17.10	---	17.30	---	17.30	---
Sigma Error (+/-)	4.82	---	3.64	---	4.83	---	3.52	---	4.14	---	3.64	---	3.84	---	4.01	---	3.95	---
MDA/LLD*	5.97	---	4.29	---	5.87	---	4.24	---	5.47	---	4.91	---	4.71	---	5.05	---	4.89	---

* Minimum Detectable Activity. Lower Level of Detection. ** Data Validation.

Parameter (pCi/g)	Table A-5.5 Gross Alpha/Beta																					
	SA5-17C-SS	DV**	SA5-17D-SS	DV**	SA5-17E-SS	DV**	SA6-18A-SS	DV**	SA6-18B-SS	DV**	SA6-18C-SS	DV**	SA6-18D-SS	DV**	SA6-18E-SS	DV**	REFC-19A-SS	DV**
Alpha	ND	---	11.40	---	9.82	---	9.77	---	8.37	---	12.90	---	ND	---	12.10	---	9.96	--
Sigma Error (+/-)	-	---	4.76	---	4.46	---	5.03	---	5.18	---	5.22	---	-	---	4.90	---	5.95	---
MDA/LLD*	8.54	u	5.29	---	5.19	---	6.57	---	7.43	---	5.90	---	6.82	u	5.33	---	8.35	---
Beta	15.20	---	16.80	---	18.00	---	16.80	---	15.10	---	12.70	---	13.90	---	14.10	---	14.80	---
Sigma Error (+/-)	3.58	---	3.39	---	3.48	---	3.43	---	3.47	---	3.13	---	3.37	---	3.10	---	3.49	---
MDA/LLD*	4.41	---	3.94	---	3.91	---	4.03	---	4.39	---	4.03	---	4.39	---	3.79	---	4.31	---

* Minimum Detectable Activity. Lower Level of Detection. ** Data Validation.

Table A-5.5 Gross Alpha/Beta								
Parameter (pCi/g)	REFC-19B-SS	DV**	REFC-19C-SS	DV**	REFC-19D-SS	DV**	REFC-19E-SS	DV**
Alpha	5.40	---	12.20	---	4.24	---	9.91	---
Sigma Error (+/-)	4.22	---	5.60	---	3.94	---	5.28	---
MDA/LLD*	6.35	U	7.21	---	6.15	U	7.12	---
Beta	11.30	---	15.60	---	9.65	---	12.60	---
Sigma Error (+/-)	3.05	---	3.44	---	2.89	---	3.27	---
MDA/LLD*	4.08	---	9.26	---	4.03	---	4.38	---

* Minimum Detectable Activity. Lower Level of Detection. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA1-1A-SS	DV**	SA1-1A-SS-D	DV**	SA1-1B-SS	DV**	SA1-1C-SS	DV**	SA1-1D-SS	DV**	SA1-1D-SS-D	DV**	SA1-1E-SS	DV**	SA1-2A-SS	DV**	SA1-2B-SS	DV**	SA1-2C-SS	DV**
Cesium 137	0.07	---	ND	---	0.06	---	0.02	---	ND	---	ND	---	0.02	---	ND	---	0.03	---	ND	---
Sigma Error (+/-)	0.02	---	-	---	0.01	---	0.01	---	-	---	-	---	0.01	---	-	---	0.01	---	-	---
MDA*	0.02	---	0.02	U	0.02	---	0.01	---	0.03	UR	0.02	U	0.02	---	0.04	U	0.01	---	0.03	U
Potassium 40	1.05	U	1.27	U	1.62	U	0.88	U	1.33	UR	1.07	U	0.96	U	ND	---	0.66	U	0.83	U
Sigma Error (+/-)	0.17	---	0.32	---	0.19	---	0.18	---	0.21	---	0.18	---	0.19	---	-	---	0.15	---	0.18	---
MDA*	0.31	---	0.97	---	0.09	---	0.24	---	0.29	---	0.22	---	0.27	---	0.49	U	0.44	---	0.40	---
Lead 212	ND	---	0.48	U	0.54	U	0.36	U	0.55	UR	0.57	U	0.59	U	0.67	U	0.40	U	ND	---
Sigma Error (+/-)	-	---	0.05	---	0.04	---	0.04	---	0.06	---	0.05	---	0.05	---	0.06	---	0.04	---	-	---
MDA*	0.03	U	0.03	---	0.03	---	0.02	---	0.03	---	0.02	---	0.02	---	0.03	---	0.02	---	0.05	U
Radium 226	0.37	U	0.41	U	0.41	U	0.25	U	0.35	UR	0.34	U	0.38	U	0.33	U	0.28	U	ND	---
Sigma Error (+/-)	0.02	---	0.04	---	0.02	---	0.02	---	0.03	---	0.02	---	0.03	---	0.03	---	0.02	---	-	---
MDA*	0.03	---	0.04	---	0.03	---	0.02	---	0.03	---	0.02	---	0.03	---	0.05	---	0.02	---	0.10	U
Radium 228	0.40	---	0.46	---	0.51	---	0.31	---	0.50	R	0.56	---	0.44	---	0.56	---	0.42	---	ND	---
Sigma Error (+/-)	0.05	---	0.06	---	0.04	---	0.03	---	0.05	---	0.05	---	0.04	---	0.05	---	0.04	---	-	---
MDA*	0.06	---	0.08	---	0.03	---	0.04	---	0.04	---	0.03	---	0.05	---	0.06	---	0.03	---	0.10	U
Thallium 208	-	---	0.15	---	0.12	---	0.16	---	0.13	R	-	-	-	-	-	-	0.11	---	-	-
Sigma Error (+/-)	-	---	0.02	---	0.07	---	0.01	---	0.05	---	-	-	-	-	-	-	0.01	---	-	-
MDA*	-	---	0.02	---	0.01	---	0.01	---	0.02	---	-	-	-	-	-	-	0.01	---	-	-
Uranium 238	ND	---	0.62	U	0.53	U	0.43	U	0.69	UR	0.51	U	0.60	U	0.67	U	0.46	U	ND	---
Sigma Error (+/-)	-	---	0.16	---	0.16	---	0.12	---	0.16	---	0.09	---	0.09	---	0.19	---	0.10	---	-	---
MDA*	0.32	U	0.30	---	0.28	---	0.18	---	0.22	---	0.15	---	0.20	---	0.34	---	0.16	---	0.52	U
Lead 210	-	-	ND	---	-	-	0.58	---	0.45	R	-	-	0.43	---	ND	---	ND	---	-	-
Sigma Error (+/-)	-	-	-	---	-	-	0.19	---	0.16	---	-	-	0.11	---	-	---	-	---	-	-
MDA*	-	-	0.39	U	-	-	0.23	---	0.32	---	-	-	0.22	---	0.63	U	0.28	U	-	-

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA1-1A-SS	DV**	SA1-1A-SS-D	DV**	SA1-1B-SS	DV**	SA1-1C-SS	DV**	SA1-1D-SS	DV**	SA1-1D-SS-D	DV**	SA1-1E-SS	DV**	SA1-2A-SS	DV**	SA1-2B-SS	DV**	SA1-2C-SS	DV**
Bismuth 212	ND	---	ND	---	0.28	---	0.28	---	0.29	R	0.42	---	ND	---	ND	---	0.29	---	ND	---
Sigma Error (+/-)	-	---	-	---	0.06	---	0.05	---	0.09	---	0.07	---	-	---	-	---	0.06	---	-	---
MDA*	0.22	U	0.12	U	0.12	---	0.10	---	0.12	---	0.08	---	0.09	U	0.18	U	0.08	---	0.26	U
Radium 224	ND	---	ND	---	0.43	---	ND	---	ND	---	0.52	---	0.30	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	0.21	---	-	---	-	---	0.29	---	0.35	---	-	---	-	---	-	---
MDA*	0.54	U	0.63	U	0.26	---	0.36	U	0.49	UR	0.18	---	0.22	---	0.69	U	0.18	U	0.71	U
Radium 223	-	-	-	-	-	-	-	-	-	-	0.17	---	-	---	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	0.04	---	-	---	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	0.06	---	-	---	-	-	-	-	-	-
Thorium 228	-	-	-	-	1.83	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	0.60	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	1.13	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	0.02	---	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	0.01	---	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	0.02	---	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.03	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.01	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.02	---
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																			
	SA1-1A-SS	DV**	SA1-1A-SS-D	DV**	SA1-1B-SS	DV**	SA1-1C-SS	DV**	SA1-1D-SS	DV**	SA1-1D-SS-D	DV**	SA1-1E-SS	DV**	SA1-2A-SS	DV**	SA1-2B-SS	DV**	SA1-2C-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																					
	SA1-2D-SS	DV**	SA1-2E-SS	DV**	SA1-3A-SS	DV**	SA1-3B-SS	DV**	SA1-3C-SS	DV**	SA1-3D-SS	DV**	SA1-3E-SS	DV**	SA1-4A-SS	DV**	SA1-4A-SS-D	DV**	SA1-4B-SS	DV**
Cesium 137	0.08	---	0.05	---	0.19	---	ND	---	0.06	---	0.21	---	0.05	---	0.35	---	0.31	---	0.18	---
Sigma Error (+/-)	0.02	---	0.01	---	0.03	---	-	---	0.01	---	0.03	---	0.01	---	0.04	---	0.03	---	0.02	---
MDA*	0.02	---	0.02	---	0.03	---	0.03	U	0.01	---	0.03	---	0.02	---	0.03	---	0.01	---	0.01	---
Potassium 40	0.51	U	0.87	U	1.08	U	0.49	U	0.85	U	1.33	U	0.60	U	3.13	---	3.09	---	1.41	---
Sigma Error (+/-)	0.09	---	0.19	---	0.18	---	0.18	---	0.18	---	0.24	---	0.12	---	0.33	---	0.18	---	0.19	---
MDA*	0.23	---	0.36	---	0.29	---	0.35	---	0.28	---	0.41	---	0.37	---	0.14	---	0.19	---	0.19	---
Lead 212	ND	---	0.50	U	0.35	U	0.46	U	0.34	U	ND	---	ND	---	0.69	---	0.70	---	0.43	---
Sigma Error (+/-)	-	---	0.05	---	0.04	---	0.04	---	0.04	---	-	---	-	---	0.06	---	0.04	---	0.04	---
MDA*	0.03	U	0.03	---	0.03	---	0.03	---	0.02	---	0.05	U	0.03	U	0.01	---	0.02	---	0.02	---
Radium 226	ND	---	0.37	U	0.31	U	0.30	U	0.29	U	0.37	U	0.29	U	0.63	U	0.74	U	0.42	U
Sigma Error (+/-)	-	---	0.03	---	0.03	---	0.03	---	0.01	---	0.03	---	0.03	---	0.04	---	0.04	---	0.02	---
MDA*	0.03	U	0.03	---	0.04	---	0.03	---	0.02	---	0.05	---	0.03	---	0.02	---	0.03	---	0.03	---
Radium 228	0.31	---	0.41	---	0.40	---	0.42	---	0.31	---	0.37	---	0.33	---	0.69	---	0.61	---	0.47	---
Sigma Error (+/-)	0.01	---	0.05	---	0.04	---	0.05	---	0.03	---	0.05	---	0.04	---	0.04	---	0.04	---	0.03	---
MDA*	0.05	---	0.06	---	0.05	---	0.09	---	0.06	---	0.08	---	0.07	---	0.03	---	0.04	---	0.05	---
Thallium 208	-	-	0.13	---	0.13	---	0.14	---	0.12	---	-	-	-	-	0.21	---	ND	---	ND	---
Sigma Error (+/-)	-	-	0.02	---	0.02	---	0.02	---	0.02	---	-	-	-	-	0.02	---	-	---	-	---
MDA*	-	-	0.02	---	0.02	---	0.02	---	0.01	---	-	-	-	-	0.01	---	0.01	U	0.01	U
Uranium 238	ND	---	0.73	U	ND	---	0.56	U	0.48	U	ND	---	0.55	U	0.96	---	0.62	---	0.56	---
Sigma Error (+/-)	-	---	0.19	---	-	---	0.14	---	0.12	---	-	---	0.13	---	0.10	---	0.09	---	0.10	---
MDA*	0.30	U	0.25	---	0.35	U	0.25	---	0.19	---	0.56	U	0.23	---	0.12	---	0.17	---	0.16	---
Lead 210	-	-	0.04	---	-	-	0.36	---	0.45	---	-	-	-	-	1.53	---	1.14	---	0.75	---
Sigma Error (+/-)	-	-	0.19	---	-	-	0.13	---	0.15	---	-	-	-	-	0.23	---	0.16	---	0.11	---
MDA*	-	-	0.34	---	-	-	0.34	---	0.24	---	-	-	-	-	0.14	---	0.22	---	0.19	---

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA1-2D-SS	DV**	SA1-2E-SS	DV**	SA1-3A-SS	DV**	SA1-3B-SS	DV**	SA1-3C-SS	DV**	SA1-3D-SS	DV**	SA1-3E-SS	DV**	SA1-4A-SS	DV**	SA1-4A-SS-D	DV**	SA1-4B-SS	DV**
Bismuth 212	0.20	---	ND	---	0.23	---	ND	---	0.22	---	ND	---	ND	---	0.40	---	0.38	---	ND	---
Sigma Error (+/-)	0.07	---	-	---	0.10	---	-	---	0.07	---	-	---	-	---	0.05	---	0.08	---	-	---
MDA*	0.11	---	0.27	U	0.14	---	0.28	U	0.10	---	0.28	U	0.25	U	0.06	---	0.10	---	0.09	U
Radium 224	ND	---	ND	---	ND	---	0.66	---	ND	---	ND	---	ND	---	-	-	-	-	-	-
Sigma Error (+/-)	-	---	-	---	-	---	0.35	---	-	---	-	---	-	---	-	-	-	-	-	-
MDA*	0.47	U	0.51	U	0.34	U	0.25	---	0.38	U	0.69	U	0.50	U	-	-	-	-	-	-
Radium 223	0.09	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	0.03	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.08	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 226	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.83	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.54	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.59	---
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.83	---	-	-	0.64	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.20	---	-	-	0.15	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.26	---	-	-	0.33	---
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																			
	SA1-2D-SS	DV**	SA1-2E-SS	DV**	SA1-3A-SS	DV**	SA1-3B-SS	DV**	SA1-3C-SS	DV**	SA1-3D-SS	DV**	SA1-3E-SS	DV**	SA1-4A-SS	DV**	SA1-4A-SS-D	DV**	SA1-4B-SS	DV**
Uranium 235	-	-		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	2.96	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.96	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	2.60	---
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA1-4C-SS	DV**	SA1-4D-SS	DV**	SA1-4E-SS	DV**	SA2-5A-SS	DV**	SA2-5A-SS-D	DV**	SA2-5B-SS	DV**	SA2-5C-SS	DV**	SA2-5D-SS	DV**	SA2-5E-SS	DV**	SA2-6A-SS	DV**
Cesium 137	ND	---	0.24	---	0.04	---	0.51	---	0.55	---	0.37	---	ND	---	0.29	---	0.06	---	0.36	---
Sigma Error (+/-)	-	---	0.02	---	0.01	---	0.05	---	0.06	---	0.03	---	-	---	0.03	---	0.02	---	0.04	---
MDA*	0.02	U	0.02	---	0.02	---	0.03	---	0.03	---	0.01	---	0.02	U	0.01	---	0.02	---	0.03	---
Potassium 40	1.94	---	1.28	---	1.41	---	2.49	U	2.71	---	3.79	---	4.62	---	2.21	U	5.51	---	2.08	U
Sigma Error (+/-)	0.28	---	0.19	---	0.14	---	0.41	---	0.32	---	0.44	---	0.60	---	0.30	---	0.65	---	0.72	---
MDA*	0.31	---	0.26	---	0.29	---	0.32	---	0.28	---	0.19	---	0.82	---	0.20	---	0.22	---	0.27	---
Lead 212	ND	---	ND	---	0.60	---	0.44	U	0.50	---	0.98	---	1.29	---	0.64	---	1.49	---	0.58	---
Sigma Error (+/-)	-	---	-	---	0.05	---	0.05	---	0.05	---	0.13	---	0.08	---	0.05	---	0.10	---	0.24	---
MDA*	0.04	U	0.02	U	0.04	---	0.05	---	0.04	---	0.02	---	0.02	---	0.02	---	0.04	---	0.03	---
Radium 226	0.44	U	ND	---	0.62	U	ND	---	0.78	U	0.84	U	1.09	---	0.64	U	1.29	---	0.76	U
Sigma Error (+/-)	0.03	---	-	---	0.04	---	-	---	0.04	---	0.04	---	0.06	---	0.04	---	0.07	---	0.04	---
MDA*	0.03	---	0.03	U	0.04	---	0.05	U	0.05	---	0.03	---	0.02	---	0.03	---	0.03	---	0.04	---
Radium 228	0.44	---	0.37	---	0.56	---	ND	---	ND	---	0.80	---	1.04	---	0.59	---	1.25	---	0.48	---
Sigma Error (+/-)	0.05	---	0.04	---	0.06	---	-	---	-	---	0.05	---	0.08	---	0.05	---	0.09	---	0.05	---
MDA*	0.08	---	0.05	---	0.07	---	0.09	U	0.07	U	0.06	---	0.03	---	0.04	---	0.07	---	0.06	---
Thallium 208	ND	---	0.10	---	0.18	---	ND	---	0.14	U	ND	---	0.32	---	0.16	---	0.42	---	0.16	---
Sigma Error (+/-)	-	---	0.01	---	0.02	---	-	---	0.02	---	-	---	0.03	---	0.02	---	0.03	---	0.02	---
MDA*	0.02	U	0.02	---	0.02	---	0.05	U	0.02	---	0.02	U	0.01	---	0.01	---	0.02	---	0.02	---
Uranium 238	ND	---	0.61	---	0.60	---	ND	---	0.72	U	0.92	U	1.54	---	0.75	U	1.63	---	0.98	U
Sigma Error (+/-)	-	---	0.09	---	0.22	---	-	---	0.20	---	0.13	---	0.20	---	0.11	---	0.22	---	0.18	---
MDA*	0.33	U	0.19	---	0.38	---	0.51	U	0.37	---	0.18	---	0.20	---	0.17	---	0.37	---	0.37	---
Lead 210	ND	---	1.25	---	ND	---	ND	---	ND	---	0.92	---	1.15	---	1.23	---	ND	---	ND	---
Sigma Error (+/-)	-	---	0.19	---	-	---	-	---	-	---	0.13	---	0.23	---	0.21	---	-	---	-	---
MDA*	2.36	U	0.19	---	2.45	U	2.78	U	2.70	U	0.23	---	0.28	---	0.24	---	1.11	U	2.50	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																					
	SA1-4C-SS	DV**	SA1-4D-SS	DV**	SA1-4E-SS	DV**	SA2-5A-SS	DV**	SA2-5A-SS-D	DV**	SA2-5B-SS	DV**	SA2-5C-SS	DV**	SA2-5D-SS	DV**	SA2-5E-SS	DV**	SA2-6A-SS	DV**
Bismuth 212	ND	---	ND	---	0.36	---	0.46	---	ND	---	0.48	---	0.72	---	0.40	---	0.75	---	ND	---
Sigma Error (+/-)	-	---	-	---	0.11	---	0.09	---	-	---	0.09	---	0.12	---	0.07	---	0.13	---	-	---
MDA*	0.13	U	0.19	U	0.12	---	0.15	---	0.28	U	0.09	---	0.10	---	0.09	---	0.17	---	0.25	U
Radium 224	-	-	-	-	-	-	ND	---	ND	---	ND	---	ND	---	-	-	ND	---	ND	---
Sigma Error (+/-)	-	-	-	-	-	-	-	---	-	---	-	---	-	---	-	-	-	---	-	---
MDA*	-	-	-	-	-	-	0.70	U	0.45	U	0.26	U	0.49	U	-	-	0.75	U	0.36	U
Radium 223	-	-	-	-	-	-	-	-	-	-	-	-	0.37	---	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	0.06	---	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	0.07	---	-	-	-	-	-	-
Thorium 228	-	-	-	-	-	-	-	-	-	-	-	-	-	-	1.06	---	3.47	---	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.55	---	0.95	---	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.31	---	1.74	---	-	-
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																					
	SA1-4C-SS	DV**	SA1-4D-SS	DV**	SA1-4E-SS	DV**	SA2-5A-SS	DV**	SA2-5A-SS-D	DV**	SA2-5B-SS	DV**	SA2-5C-SS	DV**	SA2-5D-SS	DV**	SA2-5E-SS	DV**	SA2-6A-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA2-6B-SS	DV**	SA2-6C-SS	DV**	SA2-6D-SS	DV**	SA2-6E-SS	DV**	SA3-7A-SS	DV**	SA3-7B-SS	DV**	SA3-7C-SS	DV**	SA3-7D-SS	DV**	SA3-7E-SS	DV**	SA3-8A-SS	DV**
Cesium 137	ND	---	0.05	---	0.03	---	0.14	---	0.40	---	0.26	---	0.11	---	0.12	---	0.11	---	0.04	---
Sigma Error (+/-)	-	---	0.01	---	0.01	---	0.02	---	0.05	---	0.07	---	0.02	---	0.02	---	0.02	---	0.01	---
MDA*	0.03	U	0.01	---	0.01	---	0.02	---	0.02	---	0.11	---	0.03	---	0.03	---	0.02	---	0.01	---
Potassium 40	ND	---	0.48	U	1.07	U	1.27	U	2.14	---	3.57	---	2.60	---	2.48	---	ND	---	0.33	---
Sigma Error (+/-)	-	---	0.11	---	0.15	---	0.13	---	0.30	---	0.97	---	0.37	---	0.38	---	-	---	0.11	---
MDA*	0.52	U	0.04	---	0.45	---	0.25	---	0.35	---	1.95	---	0.44	---	0.04	---	0.79	U	0.28	---
Lead 212	ND	---	0.23	U	0.36	U	0.33	U	0.57	---	ND	---	0.54	---	ND	---	0.50	---	ND	---
Sigma Error (+/-)	-	---	0.03	---	0.03	---	0.03	---	0.06	---	-	---	0.06	---	-	---	0.04	---	-	---
MDA*	0.03	U	0.02	---	0.03	---	0.03	---	0.03	---	0.19	U	0.04	---	0.04	U	0.03	---	0.02	U
Radium 226	0.16	U	0.27	U	0.33	U	0.42	U	0.54	U	ND	---	0.50	U	0.45	U	0.55	U	0.20	U
Sigma Error (+/-)	0.02	---	0.02	---	0.02	---	0.03	---	0.04	---	-	---	0.04	---	0.03	---	0.03	---	0.02	---
MDA*	0.03	---	0.02	---	0.03	---	0.02	---	0.03	---	0.19	U	0.05	---	0.04	---	0.04	---	0.02	---
Radium 228	ND	---	0.20	---	0.34	---	0.35	---	0.45	---	ND	---	ND	---	ND	---	0.51	---	ND	---
Sigma Error (+/-)	-	---	0.03	---	0.03	---	0.04	---	0.05	---	-	---	-	---	-	---	0.05	---	-	---
MDA*	0.06	U	0.04	---	0.05	---	0.05	---	0.06	---	0.31	U	0.15	U	0.09	U	0.07	---	0.05	U
Thallium 208	ND	---	0.07	U	0.09	U	0.12	U	ND	---	ND	---	0.15	---	0.14	---	ND	---	ND	---
Sigma Error (+/-)	-	---	0.01	---	0.01	---	0.01	---	-	---	-	---	0.02	---	0.03	---	-	---	-	---
MDA*	0.04	U	0.01	---	0.02	---	0.02	---	0.02	U	0.12	U	0.02	---	0.02	---	0.02	U	0.01	U
Uranium 238	ND	---	0.37	U	0.30	U	0.34	U	0.75	---	ND	---	ND	---	1.03	---	0.53	---	0.21	---
Sigma Error (+/-)	-	---	0.09	---	0.15	---	0.12	---	0.18	---	-	---	-	---	0.29	---	0.12	---	0.09	---
MDA*	0.29	U	0.14	---	0.28	---	0.31	---	0.25	---	1.18	U	0.31	U	0.38	---	0.23	---	0.19	---
Lead 210	ND	---	0.48	U	0.71	U	ND	---	1.09	---	1.18	U	ND	---	ND	---	0.48	---	ND	---
Sigma Error (+/-)	-	---	0.11	---	0.30	---	-	---	0.25	---	0.50	---	-	---	-	---	0.15	---	-	---
MDA*	1.81	U	0.19	---	0.54	---	2.19	U	0.27	---	1.16	---	0.36	U	0.61	U	0.28	---	0.31	U

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA2-6B-SS	DV**	SA2-6C-SS	DV**	SA2-6D-SS	DV**	SA2-6E-SS	DV**	SA3-7A-SS	DV**	SA3-7B-SS	DV**	SA3-7C-SS	DV**	SA3-7D-SS	DV**	SA3-7E-SS	DV**	SA3-8A-SS	DV**
Bismuth 212	ND	---	ND	---	ND	---	0.20	---	ND	---	ND	---	ND	---	ND	---	0.39	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	0.05	---	-	---	-	---	-	---	-	---	0.10	---	-	---
MDA*	0.15	U	0.13	U	0.18	U	0.15	---	0.24	U	1.12	U	0.30	U	0.36	U	0.11	---	0.16	U
Radium 224	ND	---	ND	---	ND	---	0.30	---	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	---	-	---	-	---	0.18	---	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.46	U	0.29	U	0.41	U	0.29	---	-	-	-	-	-	-	-	-	-	-	-	-
Radium 223	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 228	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																			
	SA2-6B-SS	DV**	SA2-6C-SS	DV**	SA2-6D-SS	DV**	SA2-6E-SS	DV**	SA3-7A-SS	DV**	SA3-7B-SS	DV**	SA3-7C-SS	DV**	SA3-7D-SS	DV**	SA3-7E-SS	DV**	SA3-8A-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA3-8B-SS	DV**	SA3-8C-SS	DV**	SA3-8D-SS	DV**	SA3-8E-SS	DV**	SA3-9A-SS	DV**	SA3-9B-SS	DV**	SA3-9C-SS	DV**	SA3-9D-SS	DV**	SA3-9E-SS	DV**	SA3-10A-SS	DV**
Cesium 137	0.09	---	0.06	---	0.15	---	0.14	---	<.09	U	0.27	---	0.24	---	0.40	---	0.56	---	0.61	---
Sigma Error (+/-)	0.01	---	0.01	---	0.02	---	0.02	---	-	-	0.10	---	0.07	---	0.11	---	0.13	---	0.13	---
MDA*	0.03	---	0.02	---	0.03	---	0.01	---	NR	-	NR	-	NR	-	NR	-	NR	-	NR	-
Potassium 40	1.60	---	0.35	---	5.12	---	1.13	---	5.87	---	5.19	---	3.19	---	6.08	---	4.36	---	3.16	J
Sigma Error (+/-)	0.22	---	0.12	---	0.63	---	0.20	---	1.78	---	1.84	---	1.02	---	1.52	---	1.55	---	1.33	---
MDA*	0.33	---	0.37	---	0.34	---	0.23	---	NR	-	NR	-	NR	-	NR	-	NR	-	NR	-
Lead 212	0.32	---	0.20	U	ND	---	0.71	---	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	0.03	---	0.03	---	-	---	0.06	---	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.03	---	0.04	---	0.04	U	0.02	---	-	-	-	-	-	-	-	-	-	-	-	-
Radium 226	ND	---	0.19	U	0.91	---	0.42	U	1.12	---	1.22	---	0.74	---	0.96	---	0.97	---	0.76	---
Sigma Error (+/-)	-	---	0.02	---	0.05	---	0.03	---	0.17	---	0.18	---	0.12	---	0.17	---	0.16	---	0.16	---
MDA*	0.04	U	0.04	---	0.04	---	0.02	---	NR	-	NR	-	NR	-	NR	-	NR	-	NR	-
Radium 228	0.32	---	ND	---	1.06	---	0.61	---	1.46	---	1.24	---	0.97	---	1.09	---	1.34	---	0.91	---
Sigma Error (+/-)	0.02	---	-	---	0.08	---	0.06	---	0.28	---	0.26	---	0.20	---	0.25	---	0.26	---	0.23	---
MDA*	0.08	---	0.05	U	0.09	---	0.03	---	NR	-	NR	-	NR	-	NR	-	NR	-	NR	-
Thallium 208	ND	---	0.07	---	0.36	---	0.21	---	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	---	0.01	---	0.03	---	0.02	---	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.02	U	0.02	---	0.02	---	0.01	---	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 238	ND	---	ND	---	1.31	---	ND	---	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	---	-	---	0.16	---	-	---	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.38	U	0.33	U	0.27	---	0.17	U	-	-	-	-	-	-	-	-	-	-	-	-
Lead 210	ND	---	ND	---	2.01	---	0.51	---	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	---	-	---	0.41	---	0.11	---	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	2.46	U	2.33	U	0.35	---	0.02	---	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA3-8B-SS	DV**	SA3-8C-SS	DV**	SA3-8D-SS	DV**	SA3-8E-SS	DV**	SA3-9A-SS	DV**	SA3-9B-SS	DV**	SA3-9C-SS	DV**	SA3-9D-SS	DV**	SA3-9E-SS	DV**	SA3-10A-SS	DV**
Bismuth 212	ND	---	ND	---	0.61	---	0.37	---	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	---	-	---	0.16	---	0.10	---	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.29	U	0.23	U	0.15	---	0.01	---	-	-	-	-	-	-	-	-	-	-	-	-
Radium 224	0.97	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	0.19	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.35	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Radium 223	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 228	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA3-8B-SS	DV**	SA3-8C-SS	DV**	SA3-8D-SS	DV**	SA3-8E-SS	DV**	SA3-9A-SS	DV**	SA3-9B-SS	DV**	SA3-9C-SS	DV**	SA3-9D-SS	DV**	SA3-9E-SS	DV**	SA3-10A-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA3-10B-SS	DV**	SA3-10C-SS	DV**	SA3-10D-SS	DV**	SA3-10E-SS	DV**	SA4-11A-SS	DV**	SA4-11B-SS	DV**	SA4-11C-SS	DV**	SA4-11D-SS	DV**	SA4-11E-SS	DV**	SA4-12A-SS	DV**
Cesium 137	0.48	---	0.68	---	0.58	---	0.40	---	0.11	---	0.02	---	0.17	---	0.08	---	0.02	---	ND	---
Sigma Error (+/-)	0.12	---	0.14	---	0.13	---	0.09	---	0.02	---	0.00	---	0.02	---	0.01	---	0.01	---	-	---
MDA*	NR	-	NR	-	NR	-	NR	-	0.02	---	0.01	---	0.01	---	0.02	---	0.02	---	0.03	U
Potassium 40	4.29	J	3.33	J	3.28	J	4.27	J	1.68	---	1.58	---	1.65	---	1.40	---	0.74	---	6.83	---
Sigma Error (+/-)	1.53	---	1.62	---	1.48	---	1.21	---	0.30	---	0.32	---	0.23	---	0.20	---	0.15	---	0.80	---
MDA*	NR	-	NR	-	NR	-	NR	-	0.38	---	0.57	---	0.18	---	0.25	---	0.24	---	0.24	---
Lead 212	-	-	-	-	-	-	-	-	0.40	---	0.40	---	0.43	---	0.37	---	0.32	---	1.85	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	0.04	---	0.04	---	0.01	---	0.03	---	0.02	---	0.11	---
MDA*	-	-	-	-	-	-	-	-	0.03	---	0.02	---	0.02	---	0.03	---	0.03	---	0.04	---
Radium 226	0.89	---	0.72	---	0.91	---	0.78	---	0.36	U	0.40	U	0.37	U	ND	---	0.23	U	1.02	U
Sigma Error (+/-)	0.16	---	0.14	---	0.16	---	0.13	---	0.03	---	0.03	---	0.02	---	-	---	0.02	---	0.06	---
MDA*	NR	-	NR	-	NR	-	NR	-	0.03	---	0.02	---	0.03	---	0.04	U	0.03	---	0.03	---
Radium 228	0.84	---	0.41	---	0.74	---	1.11	---	0.32	---	0.32	---	0.37	---	ND	---	0.37	---	1.75	---
Sigma Error (+/-)	0.21	---	0.21	---	0.22	---	0.22	---	0.04	---	0.03	---	0.03	---	-	---	0.04	---	0.12	---
MDA*	NR	-	NR	-	NR	-	NR	-	0.06	---	0.05	---	0.05	---	0.06	U	0.06	---	0.06	---
Thallium 208	-	-	-	-	-	-	-	-	0.12	---	0.10	---	ND	---	0.09	---	0.09	---	0.53	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	0.03	---	0.01	---	-	---	0.02	---	0.02	---	0.04	---
MDA*	-	-	-	-	-	-	-	-	0.05	---	0.01	---	0.01	U	0.01	---	0.02	---	0.02	---
Uranium 238	-	-	-	-	-	-	-	-	0.51	---	0.71	---	0.47	---	ND	---	ND	---	1.80	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	0.18	---	0.11	---	0.10	---	-	---	-	---	0.24	---
MDA*	-	-	-	-	-	-	-	-	0.30	---	0.02	---	0.15	---	0.31	U	0.31	U	0.23	---
Lead 210	-	-	-	-	-	-	-	-	1.16	---	0.50	---	0.79	---	ND	---	ND	---	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	0.19	---	0.28	---	0.13	---	-	---	-	---	-	-
MDA*	-	-	-	-	-	-	-	-	0.35	---	0.22	---	0.19	---	2.11	U	2.02	U	-	-

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA3-10B-SS	DV**	SA3-10C-SS	DV**	SA3-10D-SS	DV**	SA3-10E-SS	DV**	SA4-11A-SS	DV**	SA4-11B-SS	DV**	SA4-11C-SS	DV**	SA4-11D-SS	DV**	SA4-11E-SS	DV**	SA4-12A-SS	DV**
Bismuth 212	-	-	-	-	-	-	-	-	ND	---	0.23	---	ND	---	ND	---	ND	---	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	---	0.06	---	-	---	-	---	-	---	-	-
MDA*	-	-	-	-	-	-	-	-	0.25	U	0.09	---	0.15	U	0.21	U	0.22	U	-	-
Radium 224	-	-	-	-	-	-	-	-	0.40	---	-	-	0.29	---	0.80	---	-	-	1.23	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	0.33	---	-	-	0.19	---	0.18	---	-	-	0.48	---
MDA*	-	-	-	-	-	-	-	-	0.28	---	-	-	0.19	---	0.03	---	-	-	0.41	---
Radium 223	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 228	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	4.35	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	1.20	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	1.77	---
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.96	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.42	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.69	---
Tin 113	-	-	-	-	-	-	-	-	0.03	---	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	0.01	---	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	0.03	---	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																					
	SA3-10B-SS	DV**	SA3-10C-SS	DV**	SA3-10D-SS	DV**	SA3-10E-SS	DV**	SA4-11A-SS	DV**	SA4-11B-SS	DV**	SA4-11C-SS	DV**	SA4-11D-SS	DV**	SA4-11E-SS	DV**	SA4-12A-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	1.35	---	3.42	---	-	-	3.56	---	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	1.03	---	0.99	---	-	-	0.92	---	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	NR	-	NR	-	-	-	NR	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA4-12A-SS-D	DV**		SA4-12B-SS	DV**		SA4-12C-SS	DV**		SA4-12D-SS	DV**		SA4-12E-SS	DV**		SA4-13A-SS	DV**		SA4-13B-SS	DV**		SA4-13C-SS	DV**		SA4-13D-SS	DV**		SA4-13E-SS	DV**	
Cesium 137	ND	---		0.04	---		0.51	---		0.53	---		0.04	---		0.05	---		0.07	---		0.08	---		0.26	---		0.19	---	
Sigma Error (+/-)	-	---		0.01	---		0.05	---		0.05	---		0.01	---		0.01	---		0.01	---		0.01	---		0.03	---		0.02	---	
MDA*	0.03	U		0.02	---		0.01	---		0.02	---		0.01	---		0.02	---		0.02	---		0.02	---		0.01	---		0.02	---	
Potassium 40	7.16	---		5.92	---		1.45	U		2.12	---		1.65	U		0.98	U		1.22	U		1.21	U		2.79	---		2.50	---	
Sigma Error (+/-)	1.03	---		0.73	---		0.19	---		0.31	---		0.22	---		0.19	---		0.20	---		0.15	---		0.36	---		0.32	---	
MDA*	1.15	---		0.10	---		0.17	---		0.26	---		0.21	---		0.23	---		0.26	---		0.23	---		0.22	---		0.24	---	
Lead 212	1.86	---		1.46	---		0.54	---		ND	---		0.58	---		ND	---		ND	---		ND	---		0.56	---		ND	---	
Sigma Error (+/-)	0.11	---		0.08	---		0.04	---		-	---		0.05	---		-	---		-	---		-	---		0.04	---		0.03	---	
MDA*	0.04	---		0.04	---		0.02	---		0.03	U		0.02	---		0.03	U		0.02	U		0.02	U		0.03	---		0.03	U	
Radium 226	1.00	U		0.86	U		0.54	U		0.50	U		0.44	U		0.28	U		0.30	U		0.31	U		0.52	U		0.82	U	
Sigma Error (+/-)	0.05	---		0.05	---		0.03	---		0.03	---		0.03	---		0.02	---		0.02	---		0.02	---		0.03	---		0.04	---	
MDA*	0.04	---		0.04	---		0.02	---		0.04	---		0.02	---		0.03	---		0.03	---		0.03	---		0.03	---		0.03	---	
Radium 228	1.76	---		1.34	---		0.49	---		0.38	---		0.49	---		0.23	---		0.25	---		0.29	---		0.47	---		0.54	---	
Sigma Error (+/-)	0.11	---		0.09	---		0.04	---		0.04	---		0.04	---		0.03	---		0.03	---		0.03	---		0.04	---		0.05	---	
MDA*	0.07	---		0.03	---		0.05	---		0.07	---		0.03	---		0.01	---		0.04	---		0.05	---		0.04	---		0.05	---	
Thallium 208	0.52	---		0.43	---		ND	---		ND	---		0.15	---		0.08	U		ND	---		0.10	U		0.15	U		ND	---	
Sigma Error (+/-)	0.04	---		0.04	---		-	---		-	---		0.02	---		0.01	---		-	---		0.01	---		0.02	---		-	---	
MDA*	0.02	---		0.02	---		0.01	U		0.02	U		0.01	---		0.01	---		0.02	U		0.02	---		0.02	---		0.02	U	
Uranium 238	1.30	U		1.37	U		0.63	U		ND	---		0.67	U		ND	---		ND	---		0.48	U		0.69	U		1.13	U	
Sigma Error (+/-)	0.23	---		0.22	---		0.10	---		-	---		0.13	---		-	---		-	---		0.16	---		0.16	---		0.17	---	
MDA*	0.39	---		0.36	---		0.16	---		0.36	U		0.02	---		0.36	U		0.27	U		0.28	---		0.28	---		0.30	---	
Lead 210	1.62	---		-	-		1.72	---		-	-		0.65	U		-	-		-	-		-	-		1.44	U		1.36	---	
Sigma Error (+/-)	0.46	---		-	-		0.21	---		-	-		0.20	---		-	-		-	-		-	-		0.33	---		0.36	---	
MDA*	0.90	---		-	-		0.21	---		-	-		0.23	---		-	-		-	-		-	-		0.61	---		0.72	---	

Nuclide (pCi/g)	SA4-12A-SS-D	DV**	SA4-12B-SS	DV**	SA4-12C-SS	DV**	SA4-12D-SS	DV**	SA4-12E-SS	DV**	SA4-13A-SS	DV**	SA4-13B-SS	DV**	SA4-13C-SS	DV**	SA4-13D-SS	DV**	SA4-13E-SS	DV**
Table A-5.6 Gamma Spectroscopy																				
Bismuth 212	1.01	---	0.51	---	-	-	-	-	0.30	---	0.17	---	-	-	-	-	0.33	---	ND	---
Sigma Error (+/-)	0.13	---	0.09	---	-	-	-	-	0.07	---	0.06	---	-	-	-	-	0.07	---	-	---
MDA*	0.18	---	0.16	---	-	-	-	-	0.09	---	0.08	---	-	-	-	-	0.12	---	0.59	U
Radium 224	1.09	---	1.16	---	0.39	---	ND	---	0.47	---	ND	---	0.31	---	0.27	---	ND	---	-	-
Sigma Error (+/-)	0.52	---	0.43	---	0.24	---	-	---	0.31	---	-	---	0.18	---	0.17	---	-	---	-	-
MDA*	0.45	---	0.38	---	0.21	---	0.35	U	0.18	---	0.44	U	0.25	---	0.26	---	0.51	U	-	-
Radium 223	0.42	---	0.32	---	-	-	-	-	-	-	0.10	---	-	-	-	-	-	-	-	-
Sigma Error (+/-)	0.07	---	0.05	---	-	-	-	-	-	-	0.03	---	-	-	-	-	-	-	-	-
MDA*	0.12	---	0.11	---	-	-	-	-	-	-	0.07	---	-	-	-	-	-	-	-	-
Thorium 228	3.26	---	2.56	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	1.12	---	0.89	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.18	---	1.64	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																					
	SA4-12A-SS-D	DV**	SA4-12B-SS	DV**	SA4-12C-SS	DV**	SA4-12D-SS	DV**	SA4-12E-SS	DV**	SA4-13A-SS	DV**	SA4-13B-SS	DV**	SA4-13C-SS	DV**	SA4-13D-SS	DV**	SA4-13E-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA4-14A-SS	DV**	SA4-14A-SS-D	DV**	SA4-14B-SS	DV**	SA4-14C-SS	DV**	SA4-14D-SS	DV**	SA4-14E-SS	DV**	SA4-15A-SS	DV**	SA4-15A-SS-D	DV**	SA4-15B-SS	DV**	SA4-15C-SS	DV**
Cesium 137	0.09	---	0.15	---	ND	---	0.06	---	0.14	---	0.03	---	0.05	---	0.05	---	0.89	---	ND	---
Sigma Error (+/-)	0.02	---	0.02	---	-	---	0.01	---	0.01	---	0.01	---	0.01	---	0.01	---	0.08	---	-	---
MDA*	0.02	---	0.01	---	0.02	U	0.01	---	0.01	---	0.02	---	0.02	---	0.01	---	0.01	---	0.02	U
Potassium 40	0.93	U	1.84	U	2.02	U	0.46	U	0.45	U	0.60	U	3.46	U	2.81	U	2.90	U	0.39	U
Sigma Error (+/-)	0.16	---	0.25	---	0.27	---	0.09	---	0.09	---	0.10	---	0.42	---	0.30	---	0.33	---	0.27	---
MDA*	0.24	---	0.08	---	0.62	---	0.23	---	0.17	---	0.22	---	0.73	---	0.17	---	0.18	---	0.18	---
Lead 212	0.34	U	0.67	---	0.74	---	0.34	U	-	-	0.25	U	0.89	---	0.82	---	0.81	---	0.94	---
Sigma Error (+/-)	0.03	---	0.05	---	0.05	---	0.03	---	-	-	0.04	---	0.08	---	0.05	---	0.05	---	0.06	---
MDA*	0.03	---	0.03	---	0.03	---	0.03	---	-	-	0.03	---	0.02	---	0.02	---	0.02	---	0.02	---
Radium 226	0.25	U	0.51	U	0.48	U	0.31	U	0.28	U	0.26	U	0.77	---	0.71	---	0.74	---	0.73	---
Sigma Error (+/-)	0.03	---	0.03	---	0.03	---	0.02	---	0.02	---	0.02	---	0.05	---	0.04	---	0.04	---	0.04	---
MDA*	0.03	---	0.03	---	0.03	---	0.02	---	0.02	---	0.03	---	0.02	---	0.02	---	0.03	---	0.02	---
Radium 228	0.34	---	0.53	---	0.68	---	0.29	---	0.18	---	0.29	---	0.79	---	0.68	---	0.64	---	0.89	---
Sigma Error (+/-)	0.03	---	0.05	---	0.05	---	0.03	---	0.02	---	0.03	---	0.06	---	0.05	---	0.05	---	0.06	---
MDA*	0.06	---	0.09	---	0.04	---	0.04	---	0.04	---	0.05	---	0.03	---	0.05	---	0.05	---	0.05	---
Thallium 208	ND	---	0.18	---	0.20	---	ND	---	ND	---	0.10	U	0.23	U	0.22	U	0.23	U	0.26	---
Sigma Error (+/-)	-	---	0.02	---	0.02	---	-	---	-	---	0.01	---	0.02	---	0.02	---	0.02	---	0.02	---
MDA*	0.02	U	0.02	---	0.02	---	0.02	U	0.01	U	0.01	---	0.01	---	0.01	---	0.02	---	0.01	---
Uranium 238	ND	---	0.66	U	0.62	U	0.31	U	0.35	U	ND	---	1.01	U	0.86	U	0.91	U	1.00	U
Sigma Error (+/-)	-	---	0.16	---	0.17	---	0.11	---	0.07	---	-	---	0.17	---	0.12	---	0.16	---	0.15	---
MDA*	0.29	U	0.28	---	0.30	---	0.26	---	0.13	---	0.36	U	0.18	---	0.17	---	0.17	---	0.17	---
Lead 210	-	-	-	-	-	-	-	-	0.78	U	-	-	0.88	---	0.71	---	2.04	---	0.81	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	0.10	---	-	-	0.19	---	0.14	---	0.25	---	0.12	---
MDA*	-	-	-	-	-	-	-	-	0.15	---	-	-	0.26	---	0.22	---	0.24	---	0.21	---

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA4-14A-SS	DV**	SA4-14A-SS-D	DV**	SA4-14B-SS	DV**	SA4-14C-SS	DV**	SA4-14D-SS	DV**	SA4-14E-SS	DV**	SA4-15A-SS	DV**	SA4-15A-SS-D	DV**	SA4-15B-SS	DV**	SA4-15C-SS	DV**
Bismuth 212	-	-	0.42	---	0.40	---	-	---	0.11	---	0.21	---	0.41	---	0.44	---	0.44	---	0.53	---
Sigma Error (+/-)	-	-	0.09	---	0.09	---	-	---	0.03	---	0.08	---	0.07	---	0.07	---	0.07	---	0.08	---
MDA*	-	-	0.11	---	0.12	---	-	---	0.08	---	0.11	---	0.11	---	0.11	---	0.09	---	0.11	---
Radium 224	0.42	---	ND	---	0.31	---	ND	---	-	-	ND	---	ND	---	ND	---	-	-	ND	---
Sigma Error (+/-)	0.22	---	-	---	0.23	---	-	---	-	-	-	---	-	---	-	---	-	-	-	---
MDA*	0.26	---	0.29	U	0.31	---	0.39	U	-	-	0.43	U	0.44	U	0.24	U	-	-	0.25	U
Radium 223	-	-	-	-	-	-	0.08	---	-	-	-	-	-	-	0.22	---	0.17	---	-	-
Sigma Error (+/-)	-	-	-	-	-	-	0.04	---	-	-	-	-	-	-	0.04	---	0.03	---	-	-
MDA*	-	-	-	-	-	-	0.07	---	-	-	-	-	-	-	0.07	---	0.07	---	-	-
Thorium 228	-	-	-	-	1.27	---	-	-	-	-	-	-	-	-	2.06	---	1.65	---	-	-
Sigma Error (+/-)	-	-	-	-	0.65	---	-	-	-	-	-	-	-	-	0.08	---	0.80	---	-	-
MDA*	-	-	-	-	1.25	---	-	-	-	-	-	-	-	-	0.68	---	0.05	---	-	-
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.61	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.23	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.38	---
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	3.19	---	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	2.44	---	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	4.95	---	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																					
	SA4-14A-SS	DV**	SA4-14A-SS-D	DV**	SA4-14B-SS	DV**	SA4-14C-SS	DV**	SA4-14D-SS	DV**	SA4-14E-SS	DV**	SA4-15A-SS	DV**	SA4-15A-SS-D	DV**	SA4-15B-SS	DV**	SA4-15C-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	0.05	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	0.04	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	0.03	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA4-15D-SS	DV**	SA4-15E-SS	DV**	SA5-16A-SS	DV**	SA5-16B-SS	DV**	SA5-16C-SS	DV**	SA5-16D-SS	DV**	SA5-16E-SS	DV**	SA5-17A-SS	DV**	SA5-17B-SS	DV**	SA5-17C-SS	DV**
Cesium 137	0.22	---	0.18	---	0.07	---	0.23	---	0.04	---	0.11	---	0.24	---	0.92	---	0.21	---	0.08	---
Sigma Error (+/-)	0.03	---	0.03	---	0.01	---	0.03	---	0.01	---	0.01	---	0.03	---	0.08	---	0.04	---	0.01	---
MDA*	0.01	---	0.01	---	0.02	---	0.01	---	0.01	---	0.01	---	0.02	---	0.01	---	0.04	---	0.01	---
Potassium 40	3.74	---	3.21	U	3.43	---	1.47	U	1.36	U	0.86	U	3.03	---	4.30	---	3.88	---	3.51	---
Sigma Error (+/-)	0.44	---	0.42	---	0.42	---	0.21	---	0.20	---	0.17	---	0.34	---	0.46	---	0.71	---	0.40	---
MDA*	0.02	---	0.21	---	0.24	---	0.18	---	0.20	---	0.21	---	0.25	---	0.87	---	3.24	---	0.16	---
Lead 212	1.07	---	1.23	---	ND	---	0.33	U	0.53	---	0.45	U	0.63	---	0.72	---	0.70	---	0.63	---
Sigma Error (+/-)	0.08	---	0.09	---	-	---	0.02	---	0.05	---	0.04	---	0.04	---	0.04	---	0.08	---	0.04	---
MDA*	0.02	---	0.02	---	0.03	U	0.02	---	0.02	---	0.02	---	0.03	---	0.03	---	0.07	---	0.02	---
Radium 226	0.96	---	0.91	---	0.48	U	0.32	U	0.44	U	0.34	U	0.64	U	0.76	U	0.54	U	0.54	U
Sigma Error (+/-)	0.06	---	0.06	---	0.03	---	0.02	---	0.03	---	0.03	---	0.04	---	0.04	---	0.07	---	0.03	---
MDA*	0.02	---	0.02	---	0.04	---	0.02	---	0.02	---	0.02	---	0.04	---	0.03	---	0.08	---	0.02	---
Radium 228	0.91	---	1.10	---	0.46	---	0.28	---	0.52	---	0.33	---	0.57	---	0.57	---	ND	---	0.55	---
Sigma Error (+/-)	0.06	---	0.09	---	0.05	---	0.03	---	0.04	---	0.04	---	0.05	---	0.05	---	-	---	0.04	---
MDA*	0.03	---	0.04	---	0.06	---	0.05	---	0.03	---	0.03	---	0.06	---	0.05	---	0.17	U	0.05	---
Thallium 208	0.28	---	0.33	---	0.15	---	ND	---	0.14	U	ND	---	ND	---	0.20	---	0.17	---	0.17	---
Sigma Error (+/-)	0.03	---	0.03	---	0.02	---	-	---	0.02	---	-	---	-	---	0.02	---	0.03	---	0.02	---
MDA*	0.01	---	0.01	---	0.02	---	0.01	U	0.01	---	0.01	U	0.02	U	0.02	---	0.04	---	0.01	---
Uranium 238	1.27	U	1.33	U	0.55	U	0.38	U	0.67	U	0.58	U	ND	---	0.76	U	ND	---	0.49	U
Sigma Error (+/-)	0.23	---	0.17	---	0.15	---	0.08	---	0.12	---	0.11	---	-	---	0.17	---	-	---	0.09	---
MDA*	0.20	---	0.20	---	0.33	---	0.14	---	0.17	---	0.15	---	0.37	U	0.31	---	0.84	U	0.16	---
Lead 210	1.12	---	1.45	---	-	-	0.65	---	0.69	---	0.87	---	-	-	1.60	---	1.36	---	0.80	U
Sigma Error (+/-)	0.24	---	0.30	---	-	-	0.10	---	0.18	---	0.21	---	-	-	0.43	---	0.38	---	0.11	---
MDA*	0.27	---	0.28	---	-	-	0.17	---	0.21	---	0.17	---	-	-	0.74	---	0.69	---	0.19	---

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA4-15D-SS	DV**	SA4-15E-SS	DV**	SA5-16A-SS	DV**	SA5-16B-SS	DV**	SA5-16C-SS	DV**	SA5-16D-SS	DV**	SA5-16E-SS	DV**	SA5-17A-SS	DV**	SA5-17B-SS	DV**	SA5-17C-SS	DV**
Bismuth 212	0.62	---	0.71	---	0.29	---	-	-	0.32	---	0.16	---	-	-	0.38	---	-	-	0.35	---
Sigma Error (+/-)	0.09	---	0.08	---	0.11	---	-	-	0.06	---	0.05	---	-	-	0.09	---	-	-	0.06	---
MDA*	0.11	---	0.11	---	0.15	---	-	-	0.08	---	0.02	---	-	-	0.11	---	-	-	0.09	---
Radium 224	ND	---	-	-	0.39	---	ND	---	0.49	---	0.52	---	ND	---	ND	---	ND	---	0.30	---
Sigma Error (+/-)	-	---	-	-	0.21	---	-	---	0.39	---	0.28	---	-	---	-	---	-	---	0.22	---
MDA*	0.46	U	-	-	0.31	---	0.18	U	0.16	---	0.15	---	0.37	U	0.57	U	1.54	U	0.21	---
Radium 223	-	-	0.27	---	-	-	-	-	-	-	-	-	-	-	0.25	---	-	-	-	-
Sigma Error (+/-)	-	-	0.05	---	-	-	-	-	-	-	-	-	-	-	0.07	---	-	-	-	-
MDA*	-	-	0.07	---	-	-	-	-	-	-	-	-	-	-	0.10	---	-	-	-	-
Thorium 228	-	-	1.61	---	-	-	-	-	-	-	0.83	---	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	0.83	---	-	-	-	-	-	-	0.57	---	-	-	-	-	-	-	-	-
MDA*	-	-	0.94	---	-	-	-	-	-	-	0.66	---	-	-	-	-	-	-	-	-
Protactinium 231	-	-	-	-	0.73	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	0.20	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	0.50	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	10.40	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	2.57	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	6.20	---

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA4-15D-SS	DV**	SA4-15E-SS	DV**	SA5-16A-SS	DV**	SA5-16B-SS	DV**	SA5-16C-SS	DV**	SA5-16D-SS	DV**	SA5-16E-SS	DV**	SA5-17A-SS	DV**	SA5-17B-SS	DV**	SA5-17C-SS	DV**
Uranium 235	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA5-17D-SS	DV**	SA5-17E-SS	DV**	SA6-18A-SS	DV**	SA6-18B-SS	DV**	SA6-18C-SS	DV**	SA6-18D-SS	DV**	SA6-18E-SS	DV**	REFC-19A-SS	DV**	REFC-19B-SS	DV**
Cesium 137	0.63	---	0.45	---	0.17	---	0.10	---	0.15	---	0.17	---	0.12	---	0.35	---	0.12	---
Sigma Error (+/-)	0.06	---	0.07	---	0.02	---	0.02	---	0.02	---	0.02	---	0.02	---	0.04	---	0.02	---
MDA*	0.02	---	0.01	---	0.01	---	0.02	---	0.02	---	0.01	---	0.02	---	0.01	---	0.01	---
Potassium 40	5.82	---	4.28	---	2.78	---	2.38	U	1.82	U	1.76	U	2.27	U	2.18	U	2.04	U
Sigma Error (+/-)	0.71	---	0.49	---	0.30	---	0.27	---	0.26	---	0.22	---	0.29	---	0.25	---	0.27	---
MDA*	1.02	---	0.79	---	0.17	---	0.28	---	0.25	---	0.17	---	0.21	---	0.17	---	0.02	---
Lead 212	0.97	---	0.73	---	ND	---	ND	---	0.43	U	ND	---	ND	---	0.47	U	0.50	U
Sigma Error (+/-)	0.06	---	0.06	---	-	---	-	---	0.03	---	-	---	-	---	0.03	---	0.04	---
MDA*	0.03	---	0.02	---	0.02	U	0.03	U	0.03	---	0.02	U	0.03	U	0.02	---	0.02	---
Radium 226	0.92	U	0.72	U	0.47	U	ND	---	0.44	U	0.46	U	0.52	U	0.42	---	0.41	---
Sigma Error (+/-)	0.05	---	0.05	---	0.02	---	-	---	0.03	---	0.02	---	0.03	---	0.02	---	0.03	---
MDA*	0.03	---	0.03	---	0.02	---	0.03	U	0.03	---	0.02	---	0.03	---	0.03	---	0.02	---
Radium 228	0.82	---	0.63	---	0.41	---	0.43	---	0.33	---	0.40	---	0.41	---	0.42	---	0.48	---
Sigma Error (+/-)	0.06	---	0.05	---	0.03	---	0.04	---	0.04	---	0.03	---	0.04	---	0.04	---	0.04	---
MDA*	0.04	---	0.04	---	0.04	---	0.06	---	0.06	---	0.04	---	0.06	---	0.05	---	0.03	---
Thallium 208	0.25	---	0.20	---	ND	---	0.14	U	0.12	U	ND	---	ND	---	0.11	U	0.14	U
Sigma Error (+/-)	0.02	---	0.02	---	-	---	0.02	---	0.02	---	-	---	-	---	0.02	---	0.02	---
MDA*	0.02	---	0.01	---	0.01	U	0.02	---	0.02	---	0.01	U	0.02	U	0.01	---	0.01	---
Uranium 238	1.13	U	0.86	U	0.50	U	0.55	U	0.60	U	0.47	U	ND	---	0.52	U	0.58	U
Sigma Error (+/-)	0.21	---	0.12	---	0.07	---	0.19	---	0.23	---	0.08	---	-	---	0.10	---	0.11	---
MDA*	0.33	---	0.18	---	0.15	---	0.33	---	0.30	---	0.15	---	0.33	U	0.15	---	0.16	---
Lead 210	1.53	---	1.37	---	0.81	---	ND	---	ND	---	0.75	---	-	-	0.68	---	0.92	---
Sigma Error (+/-)	0.46	---	0.24	---	0.11	---	-	---	-	---	0.10	---	-	-	0.10	---	0.19	---
MDA*	0.77	---	0.23	---	0.19	---	2.27	U	2.13	U	0.19	---	-	-	0.19	---	0.22	---

* Minimum Detectable Activity. ** Data Validation.

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	SA5-17D-SS	DV**	SA5-17E-SS	DV**	SA6-18A-SS	DV**	SA6-18B-SS	DV**	SA6-18C-SS	DV**	SA6-18D-SS	DV**	SA6-18E-SS	DV**	REFC-19A-SS	DV**	REFC-19B-SS	DV**
Bismuth 212	0.44	---	0.42	---	0.24	---	ND	---	ND	---	-	-	0.26	---	ND	---	0.24	---
Sigma Error (+/-)	0.10	---	0.08	---	0.06	---	-	---	-	---	-	-	0.09	---	-	---	0.06	---
MDA*	0.14	---	0.09	---	0.09	---	0.21	U	0.23	U	-	-	0.11	---	0.18	U	0.10	---
Radium 224	ND	---	ND	---	ND	---	1.46	---	ND	---	ND	---	ND	---	0.26	---	-	-
Sigma Error (+/-)	-	---	-	---	-	---	0.22	---	-	---	-	---	-	---	0.18	---	-	-
MDA*	0.63	U	0.40	U	0.21	U	0.33	---	0.48	U	0.19	U	0.03	U	0.20	U	-	-
Radium 223	-	-	0.19	---	-	-	-	-	-	-	-	-	0.15	---	-	-	0.16	---
Sigma Error (+/-)	-	-	0.03	---	-	-	-	-	-	-	-	-	0.04	---	-	-	0.03	---
MDA*	-	-	0.07	---	-	-	-	-	-	-	-	-	0.09	---	-	-	0.05	---
Thorium 228	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	ND	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	---
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	0.76	U
Protactinium 231	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Nuclide (pCi/g)	Table A-5.6 Gamma Spectroscopy																	
	SA5-17D-SS	DV**	SA5-17E-SS	DV**	SA6-18A-SS	DV**	SA6-18B-SS	DV**	SA6-18C-SS	DV**	SA6-18D-SS	DV**	SA6-18E-SS	DV**	REFC-19A-SS	DV**	REFC-19B-SS	DV**
Uranium 235	-	-	0.09	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	0.03	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	0.07	---	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Bismuth 214	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Protactinium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Thorium 234	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Table A-5.6 Gamma Spectroscopy						
Nuclide (pCi/g)	REFC-19C-SS	DV**	REFC-19D-SS	DV**	REFC-19E-SS	DV**
Cesium 137	0.38	---	0.09	---	0.26	---
Sigma Error (+/-)	0.04	---	0.01	---	0.03	---
MDA*	0.01	---	0.01	---	0.01	---
Potassium 40	5.39	---	1.00	U	2.88	U
Sigma Error (+/-)	0.60	---	0.16	---	0.36	---
MDA*	0.19	---	0.25	---	0.68	---
Lead 212	ND	---	ND	---	0.53	U
Sigma Error (+/-)	-	---	-	---	0.05	---
MDA*	0.02	U	0.03	U	0.02	---
Radium 226	0.73	---	0.26	---	0.49	---
Sigma Error (+/-)	0.04	---	0.02	---	0.03	---
MDA*	0.03	---	0.03	---	0.02	---
Radium 228	0.72	---	0.25	---	0.48	---
Sigma Error (+/-)	0.05	---	0.03	---	0.04	---
MDA*	0.05	---	0.03	---	0.04	---
Thallium 208	-	-	ND	---	0.14	U
Sigma Error (+/-)	-	-	-	---	0.02	---
MDA*	-	-	0.01	U	0.01	---
Uranium 238	0.88	U	0.41	U	0.61	U
Sigma Error (+/-)	0.14	---	0.15	---	0.12	---
MDA*	0.18	---	0.23	---	0.17	---
Lead 210	2.05	---	ND	---	1.26	---
Sigma Error (+/-)	0.26	---	-	---	0.25	---
MDA*	0.24	---	0.68	U	0.23	---

Table A-5.6 Gamma Spectroscopy

Nuclide (pCi/g)	REFC-19C-SS	DV**	REFC-19D-SS	DV**	REFC-19E-SS	DV**
Bismuth 212	0.46	---	0.17	---	0.41	---
Sigma Error (+/-)	0.09	---	0.07	---	0.08	---
MDA*	0.10	---	0.09	---	0.08	---
Radium 224	0.59	---	ND	---	ND	---
Sigma Error (+/-)	0.48	---	-	---	-	---
MDA*	0.26	---	0.38	U	0.16	U
Radium 223	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-
MDA*	-	-	-	-	-	-
Thorium 228	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-
MDA*	-	-	-	-	-	-
Protactinium 231	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-
MDA*	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-
MDA*	-	-	-	-	-	-
Mercury 203	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-
MDA*	-	-	-	-	-	-
Uranium 234	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-
MDA*	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Compound (mg/Kg)	Table A-6.1 Metals/Cyanide																							
	SA1-1-V	Q*	DV**	SA1-1-V-D	Q*	DV**	SA1-2-V	Q*	DV**	SA1-3-V	Q*	DV**	SA1-4-V	Q*	DV**	SA2-5-V	Q*	DV**	SA2-6-V	Q*	DV**
Aluminum	61.6		---	71.6		---	80.9		---	129.0		---	159.0		---	183.0		-	115.0		-
Antimony	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---
Arsenic	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ
Barium	2.8	B	J	7.6	B	J	14.9	B	J	10.6	B	J	16.5	B	J	6.4	B	J	7.1	B	J
Beryllium	0.19	B	U	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---
Cadmium	0.83	B	U	0.83	B	U	1.10		U	1.10		U	0.62	U	---	0.62	U	---	0.62	U	---
Calcium	2430		---	2240		---	1610		---	1760		---	2210		---	2260		---	923	B	J
Chromium	0.92	U	---	0.92	U	---	0.92	U	---	0.92	U	---	0.92	U	---	5.90		---	0.92	U	---
Cobalt	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---
Copper	3.4	B	U	3.3	B	U	2.8	B	U	2.2	B	U	2.8	B	U	6.7		---	3.0	B	U
Iron	37.3		---	19.8	B	J	24.1		---	16	B	J	25.3		---	48.7		---	13.9	B	J
Lead	2.00		---	0.36	B	U	1.30		---	2.60		---	3.70	S	---	1.60		---	3.30		---
Magnesium	320	B	J	284	B	J	308	B	J	294	B	J	328	B	J	385	B	J	238	B	J
Manganese	63.7		---	146.0		---	155.0		---	145.0		---	140.0		---	194.0		---	217.0		---
Mercury	0.05	U	UJ	0.05	U	---	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ
Nickel	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---
Potassium	1090		---	1180		---	762	B	J	597	U	---	676	B	J	910	B	J	647	B	J
Selenium	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---
Silver	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---
Sodium	48.8	B	J	68.2	B	J	45.7	B	J	42.7	B	J	65.1	B	J	48.4	B	J	41.6	B	J
Thallium	0.20	UW	UJ	0.20	UW	J	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ
Vanadium	0.42	U	---	0.42	U	---	0.42	U	---	0.42	U	---	0.42	U	---	0.81	B	U	0.84	B	U
Zinc	22.3	E	J	18.5	E	J	20.4	E	J	29.5	E	J	13.2	E	J	19.5	E	J	27.1	E	J
Cyanide	0.13	U	UJ	0.13	U	UJ	0.13	U	UJ	0.12	U	UJ	0.13	U	UJ	0.13	U	UJ	0.13	U	UJ

* Data Qualifier. ** Data Validation.

Table A-6.1 Metals/Cyanide

Compound (mg/Kg)	SA3-7-V	Q*	DV**	SA3-8-V	Q*	DV**	SA3-9-V	Q*	DV**	SA3-10-V	Q*	DV**	SA4-11-V	Q*	DV**	SA4-12-V	Q*	DV**	SA4-13-V	Q*	DV**
Aluminum	99.9		---	115.0		---	87.6		---	97.7		---	47.6		---	108.0		---	117.0		---
Antimony	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---
Arsenic	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ
Barium	6.2	B	J	7.6	B	J	7.5	B	J	6.7	B	J	3.4	B	J	7.3	B	J	3.9	B	J
Beryllium	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---
Cadmium	0.62	U	---	0.62	U	---	0.62	U	---	0.75	B	U	0.70	B	U	0.62	U	---	0.83	B	U
Calcium	1440		---	2050		---	2270		---	1260		---	2180		---	1470		---	3280		---
Chromium	0.92	U	---	0.92	U	---	0.92	U	---	0.92	U	---	0.92	U	---	0.92	U	---	0.92	U	---
Cobalt	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---
Copper	2.8	B	U	2.6	B	U	3.2	B	U	2.6	B	U	2.6	B	U	2.6	B	U	3.6	B	U
Iron	23.6		---	20.9		---	25.8		---	16.3	B	J	22.7		---	13.2	B	J	29.8		---
Lead	14.30		---	2.90	S	---	1.60		---	13.00		---	4.30	S	---	0.57	B	UJ	0.48	B	U
Magnesium	240	B	J	251	B	J	281	B	J	225	B	J	277	B	J	217	B	J	341	B	J
Manganese	157.0		---	168.0		---	87.9		---	115.0		---	33.4		---	115.0		---	88.1		---
Mercury	0.05	U	---	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ
Nickel	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---
Potassium	597	U	---	597	U	---	606	B	J	754	B	J	1000	B	J	623	B	J	688	B	J
Selenium	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---	4.0	U	---
Silver	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---
Sodium	61.3	B	J	60.0	B	J	53.2	B	J	37.1	B	J	53.2	B	J	50.8	B	J	63.7	B	J
Thallium	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	U	---	0.20	U	---
Vanadium	0.42	U	---	0.83	B	U	0.42	U	---	0.43	B	U	0.42	U	---	0.42	U	---	0.42	U	---
Zinc	16.5	E	J	11.8	E	J	11.7	E	J	8.4	E	J	12.5	E	J	8.6	E	J	17.5	E	J
Cyanide	0.13	U	UJ	0.13	U	UJ	0.12	U	UJ	0.13	U	UJ	0.13	U	UJ	0.13	U	UJ	0.13	U	UJ

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	Table A-6.1 Metals/Cyanide																	
	SA4-14-V	Q*	DV**	SA4-15-V	Q*	DV**	SA5-16-V	Q*	DV**	SA5-17-V	Q*	DV**	SA6-18-V	Q*	DV**	REFC-19-V	Q*	DV**
Aluminum	97.0		---	131.0		---	159.0		---	244.0		---	238.0		---	207.0		---
Antimony	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---	11.7	U	---
Arsenic	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ
Barium	9.0	B	J	4.6	B	J	8.0	B	J	17.2	B	J	169.0		---	8.0	B	J
Beryllium	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---	0.06	U	---
Cadmium	0.62	U	---	0.62	U	---	0.62	U	---	0.96	B	U	0.62	U	---	0.65	B	U
Calcium	1750		---	2070		---	2330		---	2700		---	4840		---	1650		---
Chromium	0.92	U	---	0.92	U	---	0.92	U	---	0.92	U	---	0.99	B	J	0.92	U	---
Cobalt	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---	0.84	U	---
Copper	3.2	B	U	5.3		U	2.8	B	U	2.8	B	U	6.3		---	3.2	B	J
Iron	14.3	B	J	21.2		---	20.7		---	28.1		---	165		---	28.9		---
Lead	0.20	B	U	0.43	B	U	0.28	B	U	1.90		---	2.70		---	0.29	B	U
Magnesium	259	B	J	227	B	J	352	B	J	394	B	J	770	B	J	289	B	J
Manganese	115.0		---	115.0		---	102.0		---	178.0		---	300.0		---	58.9		---
Mercury	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ	0.05	U	UJ
Nickel	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---	4.2	U	---
Potassium	659	B	J	877	B	J	701	B	J	820	B	J	623	B	J	1010		---
Selenium	4.0	U	---	4.0	U	---	4.0	U	---	4.0	UW	UJ	4.0	U	---	4.0	UW	UJ
Silver	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---	0.70	U	---
Sodium	36.1	B	J	64.2	B	J	48.5	B	J	52.9	B	J	52.5	B	J	38.2	B	J
Thallium	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ	0.20	UW	UJ
Vanadium	0.42	U	---	0.42	U	---	0.42	U	---	0.42	U	---	0.76	B	U	0.42	U	---
Zinc	42.0	E	J	115.0	E	J	102.0	E	J	11.0	E	J	43.5	E	J	12.2	E	J
Cyanide	0.13	U	UJ	0.13	U	UJ	0.13	U	UJ	0.13	U	UJ	0.13	U	UJ	0.11	U	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-6.2 Pesticides/PCB's																							
	SA1-1-V	Q*	DV**	SA1-1-V-D	Q*	DV**	SA1-2-V	Q*	DV**	SA1-3-V	Q*	DV**	SA1-4-V	Q*	DV**	SA2-5-V	Q*	DV**	SA2-6-V	Q*	DV**
Alpha-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Beta-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Delta-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Gamma-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Heptachlor	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Aldrin	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Heptachlor epoxide	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Endosulfan I	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Dieldrin	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
4,4'-DDE	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Endrin	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Endosulfan II	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
4,4'-DDD	32	U	UJ	32	U	UJ	50	X	J	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Endosulfan sulfate	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	42		J	32	U	UJ
4,4'-DDT	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Methoxychlor	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Endrin ketone	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Alpha-chlordane	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Gamma-chlordane	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Toxaphene	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ
Aroclor-1016	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1221	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1232	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1242	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	SA1-1-V	Q*	DV**	SA1-1-V-D	Q*	DV**	SA1-2-V	Q*	DV**	SA1-3-V	Q*	DV**	SA1-4-V	Q*	DV**	SA2-5-V	Q*	DV**	SA2-6-V	Q*	DV**
Table A-6.2 Pesticides/PCB's																					
Aroclor-1248	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1254	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ
Aroclor-1260	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ

Compound (ug/Kg)	Table A-6.2 Pesticides/PCB's																							
	SA3-7-V	Q*	DV**	SA3-8-V	Q*	DV**	SA3-9-V	Q*	DV**	SA3-10-V	Q*	DV**	SA4-11-V	Q*	DV**	SA4-12-V	Q*	DV**	SA4-13-V	Q*	DV**
Alpha-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Beta-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	27		J
Delta-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Gamma-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Heptachlor	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Aldrin	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Heptachlor epoxide	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Endosulfan I	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ
Dieldrin	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
4,4'-DDE	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Endrin	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Endosulfan II	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
4,4'-DDD	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Endosulfan sulfate	32	U	UJ	35		J	39		J	42		J	39		J	45		J	32	U	UJ
4,4'-DDT	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ
Methoxychlor	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Endrin ketone	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	70		J	70		J	130		J
Alpha-chlordane	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Gamma-chlordane	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Toxaphene	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ
Aroclor-1016	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1221	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1232	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1242	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	SA3-7-V	Q*	DV**	SA3-8-V	Q*	DV**	SA3-9-V	Q*	DV**	SA3-10-V	Q*	DV**	SA4-11-V	Q*	DV**	SA4-12-V	Q*	DV**	SA4-13-V	Q*	DV**
	Table A-6.2 Pesticides/PCB's																				
Aroclor-1248	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ
Aroclor-1254	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ
Aroclor-1260	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-6.2 Pesticides/PCB's																				
	SA4-14-V	Q*	DV**	SA4-15-V	Q*	DV**	SA5-16-V	Q*	DV**	SA5-17-V	Q*	DV**	SA6-18-V	Q*	DV**	REFC-19-V	Q*	DV**			
Alpha-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Beta-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Delta-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Gamma-BHC	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Heptachlor	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Aldrin	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Heptachlor epoxide	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Endosulfan I	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ	16	U	UJ			
Dieldrin	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ			
4,4'-DDE	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ			
Endrin	40		J	52		J	42		J	32	U	UJ	32	U	UJ	32	U	UJ			
Endosulfan II	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ			
4,4'-DDD	32	U	UJ	32	U	UJ	32	U	UJ	88	X	J	32	U	UJ	32	U	UJ			
Endosulfan sulfate	32	U	UJ	71		J	49		J	32	U	UJ	32	U	UJ	32	U	UJ			
4,4'-DDT	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ	32	U	UJ			
Methoxychlor	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			
Endrin ketone	32	U	UJ	88		J	87		J	87		J	32	U	UJ	32	U	UJ			
Alpha-chlordane	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			
Gamma-chlordane	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			
Toxaphene	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ			
Aroclor-1016	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			
Aroclor-1221	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			
Aroclor-1232	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			
Aroclor-1242	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-6.2 Pesticides/PCB's																				
	SA4-14-V	Q*	DV**	SA4-15-V	Q*	DV**	SA5-16-V	Q*	DV**	SA5-17-V	Q*	DV**	SA6-18-V	Q*	DV**	REFC-19-V	Q*	DV**			
Aroclor-1248	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ	160	U	UJ			
Aroclor-1254	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ			
Aroclor-1260	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ	320	U	UJ			

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	SA1-1-V	Q*	DV**	SA1-1-V-D	Q*	DV**	SA1-2-V	Q*	DV**	SA1-3-V	Q*	DV**	SA1-4-V	Q*	DV**	SA2-5-V	Q*	DV**	SA2-6-V	Q*	DV**
	Table A-6.3 Semi-Volatiles																				
Phenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
bis (2-Cloroethyl)Ether	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Chlorophenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,3-Dichlorobenzene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,4-Dichlorobenzene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzyl Alcohol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,2-Dichlorobenzene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Methylphenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
bis (2-Chloroisopropyl)Ether	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Methylphenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
N-Nitroso-Di-n-Propylamine	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachloroethane	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Nitrobenzene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Isophorone	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Nitrophenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dimethylphenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzoic Acid	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
bis (2-Chloroethoxy)Methane	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dichlorophenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,2,4-Trichlorobenzene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Naphthalene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Chloroaniline	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachlorobutadiene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Chloro-3-Methylphenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ

Table A-6.3 Semi-Volatiles

Compound (ug/Kg)	SA1-1-V	Q*	DV**	SA1-1-V-D	Q*	DV**	SA1-2-V	Q*	DV**	SA1-3-V	Q*	DV**	SA1-4-V	Q*	DV**	SA2-5-V	Q*	DV**	SA2-6-V	Q*	DV**
2-Methylnaphthalene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachlorocyclopentadiene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4,6-Trichlorophenol	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4,5-Trichlorophenol	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
2-Chloronaphthalene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Nitroaniline	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
Dimethyl Phthalate	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Acenaphthylene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,6-Dinitrotoluene	6600	U	J	1300	J	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	960	J	UJ	6600	U	UJ
3-Nitroaniline	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
Acenaphthene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dinitrophenol	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
4-Nitrophenol	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
Dibenzofuran	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dinitrotoluene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Diethylphthalate	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Chlorophenyl-phenylether	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Fluorene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Nitroaniline	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
4,6-Dinitro-2-Methylphenol	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
N-Nitrosodiphenylamine (1)	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Bromophenyl-phenylether	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachlorobenzene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Pentachlorophenol	32000	U	J	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ

* Data Qualifier. ** Data Validation.

Table A-6.3 Semi-Volatiles

Compound (ug/Kg)	SA1-1-V	Q*	DV**	SA1-1-V-D	Q*	DV**	SA1-2-V	Q*	DV**	SA1-3-V	Q*	DV**	SA1-4-V	Q*	DV**	SA2-5-V	Q*	DV**	SA2-6-V	Q*	DV**
Phenanthrene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Anthracene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Di-n-Butylphthalate	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Fluoranthene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Pyrene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Butylbenzylphthalate	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
3,3'-Dichlorbenzidine	13000	U	J	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ
Benzo(a)Anthracene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Chrysene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Bis (2-Ethylhexyl)Phthalate	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Di-n-Octyl Phthalate	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(b)Fluoranthene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(k)Fluoranthene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(a)Pyrene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Indeno(1,2,3-cd)Pyrene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Dibenz(a,h)Anthracene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(g,h,i)Perylene	6600	U	J	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ

Compound (ug/Kg)	Table A-6.3 Semi-Volatiles																							
	SA3-7-V	Q*	DV**	SA3-8-V	Q*	DV**	SA3-9-V	Q*	DV**	SA3-10-V	Q*	DV**	SA4-11-V	Q*	DV**	SA4-12-V	Q*	DV**	SA4-13-V	Q*	DV**
Phenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	3100	J	---	6600	U	UJ	6600	U	UJ
bis (2-Cloroethyl)Ether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Chlorophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,3-Dichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,4-Dichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzyl Alcohol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,2-Dichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Methylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
bis (2-Chloroisopropyl)Ether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Methylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
N-Nitroso-Di-n-Propylamine	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachloroethane	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Nitrobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Isophorone	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Nitrophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dimethylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzoic Acid	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	14000	J	---	9900	J	---	32000	U	UJ
bis (2-Chloroethoxy)Methane	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dichlorophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
1,2,4-Trichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Naphthalene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Chloroaniline	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachlorobutadiene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Chloro-3-Methylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ

Compound (ug/Kg)	Table A-6.3 Semi-Volatiles																							
	SA3-7-V	Q*	DV**	SA3-8-V	Q*	DV**	SA3-9-V	Q*	DV**	SA3-10-V	Q*	DV**	SA4-11-V	Q*	DV**	SA4-12-V	Q*	DV**	SA4-13-V	Q*	DV**
2-Methylnaphthalene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachlorocyclopentadiene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4,6-Trichlorophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4,5-Trichlorophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
2-Chloronaphthalene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2-Nitroaniline	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
Dimethyl Phthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Acenaphthylene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,6-Dinitrotoluene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
3-Nitroaniline	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
Acenaphthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dinitrophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
4-Nitrophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
Dibenzofuran	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
2,4-Dinitrotoluene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Diethylphthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Chlorophenyl-phenylether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Fluorene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Nitroaniline	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
4,6-Dinitro-2-Methylphenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ
N-Nitrosodiphenylamine (1)	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
4-Bromophenyl-phenylether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Hexachlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Pentachlorophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-6.3 Semi-Volatiles																							
	SA3-7-V	Q*	DV**	SA3-8-V	Q*	DV**	SA3-9-V	Q*	DV**	SA3-10-V	Q*	DV**	SA4-11-V	Q*	DV**	SA4-12-V	Q*	DV**	SA4-13-V	Q*	DV**
Phenanthrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Anthracene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Di-n-Butylphthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Fluoranthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Pyrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Butylbenzylphthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
3,3'-Dichlorbenzidine	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ
Benzo(a)Anthracene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Chrysene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Bis (2-Ethylhexyl)Phthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ
Di-n-Octyl Phthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	R	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(b)Fluoranthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	R	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(k)Fluoranthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	R	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(a)Pyrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	R	6600	U	UJ	6600	U	UJ	6600	U	UJ
Indeno(1,2,3-cd)Pyrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	R	6600	U	UJ	6600	U	UJ	6600	U	UJ
Dibenz(a,h)Anthracene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	R	6600	U	UJ	6600	U	UJ	6600	U	UJ
Benzo(g,h,i)Perylene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	R	6600	U	UJ	6600	U	UJ	6600	U	UJ

Table A-6.3 Semi-Volatiles

Compound (ug/Kg)	SA4-14-V	Q*	DV**	SA4-15-V	Q*	DV**	SA5-16-V	Q*	DV**	SA5-17-V	Q*	DV**	SA6-18-V	Q*	DV**	REFC-19-V	Q*	DV**
Phenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
bis (2-Cloroethyl)Ether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2-Chlorophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
1,3-Dichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
1,4-Dichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Benzyl Alcohol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
1,2-Dichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2-Methylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
bis (2-Chloroisopropyl)Ether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
4-Methylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
N-Nitroso-Di-n-Propylamine	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Hexachloroethane	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Nitrobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Isophorone	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2-Nitrophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2,4-Dimethylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Benzoic Acid	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
bis (2-Chloroethoxy)Methane	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2,4-Dichlorophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
1,2,4-Trichlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Naphthalene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
4-Chloroaniline	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Hexachlorobutadiene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
4-Chloro-3-Methylphenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-6.3 Semi-Volatiles																				
	SA4-14-V	Q*	DV**	SA4-15-V	Q*	DV**	SA5-16-V	Q*	DV**	SA5-17-V	Q*	DV**	SA6-18-V	Q*	DV**	REFC-19-V	Q*	DV**
2-Methylnaphthalene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Hexachlorocyclopentadiene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2,4,6-Trichlorophenol	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2,4,5-Trichlorophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
2-Chloronaphthalene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2-Nitroaniline	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
Dimethyl Phthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Acenaphthylene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2,6-Dinitrotoluene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
3-Nitroaniline	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
Acenaphthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2,4-Dinitrophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
4-Nitrophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
Dibenzofuran	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
2,4-Dinitrotoluene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Diethylphthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
4-Chlorophenyl-phenylether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Fluorene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
4-Nitroaniline	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
4,6-Dinitro-2-Methylphenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J
N-Nitrosodiphenylamine (1)	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
4-Bromophenyl-phenylether	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Hexachlorobenzene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Pentachlorophenol	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	UJ	32000	U	J

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-6.3 Semi-Volatiles																				
	SA4-14-V	Q*	DV**	SA4-15-V	Q*	DV**	SA5-16-V	Q*	DV**	SA5-17-V	Q*	DV**	SA6-18-V	Q*	DV**	REFC-19-V	Q*	DV**
Phenanthrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Anthracene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Di-n-Butylphthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Fluoranthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Pyrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Butylbenzylphthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
3,3'-Dichlorbenzidine	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	UJ	13000	U	J
Benzo(a)Anthracene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Chrysene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Bis (2-Ethylhexyl)Phthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Di-n-Octyl Phthalate	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Benzo(b)Fluoranthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Benzo(k)Fluoranthene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Benzo(a)Pyrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Indeno(1,2,3-cd)Pyrene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Dibenz(a,h)Anthracene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J
Benzo(g,h,i)Perylene	6600	U	UJ	6600	U	UJ	6600	U	UJ	6600	U	UJ	6500	U	UJ	6600	U	J

Table A-6.4 Gross Alpha/Beta

Parameter (pCi/g)	SA1-1-V	DV**	SA1-1-V-D	DV**	SA1-2-V	DV**	SA1-3-V	DV**	SA1-4-V	DV**	SA2-5-V	DV**	SA2-6-V	DV**	SA3-7-V	DV**	SA3-8-V	DV**
Alpha	0.33	---	0.29	---	0.42	---	ND	---	0.29	---	0.40	---	0.26	---	0.28	---	0.42	---
Sigma Error (+/-)	1.22	---	0.34	---	0.32	---	-	---	0.39	---	3.50	---	0.36	---	0.25	---	1.64	---
MDA*	0.24	---	0.26	---	0.26	---	0.24	U	0.23	---	0.21	---	0.21	---	0.24	---	0.21	---
Beta	0.41	---	0.22	---	0.47	---	0.54	---	0.28	---	0.54	---	0.31	---	0.40	---	0.37	---
Sigma Error (+/-)	0.06	---	0.03	---	0.07	---	0.08	---	0.04	---	0.08	---	0.05	---	0.06	---	0.06	---
MDA*	0.20	---	0.20	---	0.20	---	0.19	---	0.20	---	0.19	---	0.18	---	0.19	---	0.18	---

* Minimum Detectable Activity. ** Data Validation.

Table A-6.4 Gross Alpha/Beta

Parameter (pCi/g)	SA3-9-V	DV**	SA3-10-V	DV**	SA4-11-V	DV**	SA4-12-V	DV**	SA4-13-V	DV**	SA4-14-V	DV**	SA4-15-V	DV**	SA5-16-V	DV**	SA5-17-V	DV**
Alpha	0.28	---	0.36	---	ND	---	0.39	---	ND	---	3.89	---	0.36	---	0.36	---	ND	---
Sigma Error (+/-)	0.49	---	0.56	---	-	---	0.44	---	-	---	1.35	---	0.27	---	0.72	---	-	---
MDA*	0.24	---	0.23	---	0.31	U	0.29	---	0.42	U	0.46	---	0.31	---	0.16	---	0.33	U
Beta	ND	---	0.64	---	0.37	---	0.49	---	ND	---	0.89	---	0.65	---	0.30	---	0.45	---
Sigma Error (+/-)	-	---	0.11	---	0.05	---	0.07	---	-	---	0.17	---	0.12	---	0.04	---	0.06	---
MDA*	0.25	U	0.21	---	0.21	---	0.25	---	0.23	U	0.22	---	0.21	---	0.19	---	0.21	---

Table A-6.4 Gross Alpha/Beta				
Parameter (pCi/g)	SA6-18-V	DV**	REFC-19-V	DV**
Alpha	2.06	---	0.28	---
Sigma Error (+/-)	2.03	---	0.32	---
MDA*	0.58	---	0.25	---
Beta	2.56	---	1.04	---
Sigma Error (+/-)	0.59	---	0.19	---
MDA*	0.40	---	0.21	---

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-6.5 Gamma Spectroscopy																	
	SA1-1-V	DV**	SA1-1-V-D	DV**	SA1-2-V	DV**	SA1-3-V	DV**	SA1-4-V	DV**	SA2-5-V	DV**	SA2-6-V	DV**	SA3-7-V	DV**	SA3-8-V	DV**
Cesium 137	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.21	U	0.22	U	0.20	U	0.27	U	0.14	U	0.19	U	0.21	U	0.21	U	0.19	U
Potassium 40	ND	---	15.40	---	ND	---	6.26	---	6.44	---	8.58	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	1.15	---	-	---	1.96	---	1.08	---	1.31	---	-	---	-	---	-	---
MDA*	3.62	U	-	---	3.09	U	4.06	---	1.41	---	1.33	---	3.12	U	3.25	U	2.79	U
Lead 212	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.32	U	0.33	U	0.22	U	0.37	U	0.23	U	0.19	U	0.21	U	0.28	U	0.24	U
Thallium 208	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.19	U	0.22	U	0.17	U	0.23	U	0.18	U	0.17	U	0.18	U	0.21	U	0.16	U
Uranium 238	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.93	U	2.18	U	1.77	U	2.26	U	1.49	U	1.56	U	1.66	U	2.56	U	2.27	U
Lead 210	2.29	---	ND	---	ND	---	3.52	---	-	-	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	0.80	---	-	---	-	---	1.09	---	-	-	-	---	-	---	-	---	-	---
MDA*	1.81	---	2.16	U	1.82	U	2.24	---	-	-	1.54	U	1.63	U	20.40	U	18.10	U
Beryllium 7	3.16	---	4.35	---	-	-	3.47	---	3.00	---	2.36	---	-	-	1.92	---	-	-
Sigma Error (+/-)	0.86	---	0.97	---	-	-	0.89	---	0.78	---	0.39	---	-	-	0.56	---	-	-
MDA*	1.19	---	-	---	-	-	1.39	---	0.80	---	0.89	---	-	-	1.19	---	-	-
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.36	U	0.51	U	0.48	U	0.50	U	0.36	U	0.41	U	0.41	U	0.39	U	0.29	U
Cerium 139	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	SA1-1-V	DV**	SA1-1-V-D	DV**	SA1-2-V	DV**	SA1-3-V	DV**	SA1-4-V	DV**	SA2-5-V	DV**	SA2-6-V	DV**	SA3-7-V	DV**	SA3-8-V	DV**
Table A-6.5 Gamma Spectroscopy																		
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Nuclide (pCi/g)	Table A-6.5 Gamma Spectroscopy																			
	SA3-9-V	DV**	SA3-10-V	DV**	SA4-11-V	DV**	SA4-12-V	DV**	SA4-13-V	DV**	SA4-14-V	DV**	SA4-15-V	DV**	SA5-16-V	DV**	SA5-17-V	DV**
Cesium 137	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	0.21	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.10	---	-	---
MDA*	0.19	U	0.14	U	0.20	U	0.21	U	0.22	U	0.31	U	0.23	U	0.18	---	0.15	U
Potassium 40	ND	---	ND	---	ND	---	ND	---	ND	---	21.30	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	3.12	---	-	---	-	---	-	---
MDA*	2.72	U	5.56	U	2.83	U	3.16	U	3.96	U	8.23	---	4.25	U	3.13	U	2.48	U
Lead 212	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.18	U	0.26	U	0.21	U	0.24	U	0.35	U	0.46	U	0.38	U	0.25	U	0.17	U
Thallium 208	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.14	U	0.14	U	0.19	U	0.22	U	0.22	U	0.30	U	0.30	U	0.21	U	0.14	U
Uranium 238	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.50	U	1.47	U	1.60	U	1.84	U	2.69	U	3.02	U	2.32	U	2.44	U	1.35	U
Lead 210	1.46	---	ND	---	ND	---	2.21	---	ND	---	ND	---	4.15	---	ND	---	2.10	---
Sigma Error (+/-)	0.44	---	-	---	-	---	0.91	---	-	---	-	---	1.05	---	-	---	0.51	---
MDA*	1.50	---	1.43	U	1.54	U	1.84	---	24.30	U	2.99	U	2.15	---	19.00	U	1.33	---
Beryllium 7	-	-	-	-	-	-	-	-	-	-	5.78	---	-	-	2.83	--	3.57	---
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	1.28	---	-	-	0.62	---	0.72	---
MDA*	-	-	-	-	-	-	-	-	-	-	1.68	---	-	-	1.00	---	0.78	---
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.44	U	0.33	U	0.33	U	0.46	U	0.47	U	0.71	U	0.53	U	0.30	U	0.29	U
Cerium 139	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-6.5 Gamma Spectroscopy																			
	SA3-9-V	DV**	SA3-10-V	DV**	SA4-11-V	DV**	SA4-12-V	DV**	SA4-13-V	DV**	SA4-14-V	DV**	SA4-15-V	DV**	SA5-16-V	DV**	SA5-17-V	DV**
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Tin 113	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
Sigma Error (+/-)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
MDA*	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-6.5 Gamma Spectroscopy

Nuclide (pCi/g)	SA6-18-V	DV**	REFC-19-V	DV**
Cesium 137	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---
MDA*	0.27	U	0.20	U
Potassium 40	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---
MDA*	5.22	U	3.71	U
Lead 212	0.53	---	ND	---
Sigma Error (+/-)	0.23	---	-	---
MDA*	0.33	---	0.26	U
Thallium 208	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---
MDA*	0.23	U	0.18	U
Uranium 238	ND	---	ND	---
Sigma Error (+/-)	-	---	-	---
MDA*	2.09	U	1.56	U
Lead 210	7.61	---	2.33	---
Sigma Error (+/-)	1.19	---	1.04	---
MDA*	2.03	---	1.47	---
Beryllium 7	25.50	---	-	-
Sigma Error (+/-)	2.78	---	-	-
MDA*	1.25	---	-	-
Bismuth 214	0.68	---	ND	---
Sigma Error (+/-)	0.23	---	-	---
MDA*	0.49	---	0.33	U
Cerium 139	0.15	---	-	-

* Minimum Detectable Activity. ** Data Validation.

Table A-6.5 Gamma Spectroscopy				
Nuclide (pCi/g)	SA6-18-V	DV**	REFC-19-V	DV**
Sigma Error (+/-)	0.09	---	-	-
MDA*	0.12	---	-	-
Tin 113	-	-	0.20	---
Sigma Error (+/-)	-	-	0.09	---
MDA*	-	-	0.18	---

* Minimum Detectable Activity. ** Data Validation.

Compound (mg/Kg)	Table A-7.1 Metals																				
	2A-SM-B-I	Q*	DV**	2A-SM-F-I	Q*	DV**	2A-SM-L-I	Q*	DV**	2A-SM-M-I	Q*	DV**	3A-SM-B-I	Q*	DV**	3A-SM-M-I	Q*	DV**	4H-SM-C-I	Q*	DV**
Aluminum	5.90	B	J	5.10	B	J	1.00	B	J	1.80	B	J	7.60	B	J	2.80	B	J	10.70		---
Antimony	2.6	U	---	2.5	U	---	2.4	U	---	2.5	U	---	2.3	U	---	2.5	U	---	2.2	U	---
Arsenic	0.42	UW	UJ	0.40	U	---	0.41	U	---	0.37	U	---	0.46	UW	UJ	0.47	U	---	0.43	U	---
Barium	56.90		---	0.30	B	J	0.14	B	J	0.16	B	J	32.10		---	0.30	B	J	14.90		---
Beryllium	0.05	B	U	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.04	B	U
Cadmium	0.13	U	---	0.12	U	---	0.12	U	---	0.12	U	---	0.11	U	---	0.12	U	---	0.11	U	---
Calcium	98000		---	251		---	139	B	J	158	B	J	98500		---	291		---	7960		---
Chromium	1.40		---	0.39	B	U	0.15	U	---	0.35	B	J	3.40		---	0.36	B	J	0.31	B	U
Cobalt	0.28	U	---	0.27	U	---	0.26	U	---	0.27	U	---	0.24	U	---	0.27	U	---	0.24	U	---
Copper	11.10		---	5.30		---	6.80		---	1.90		---	0.81	B	U	2.20		---	3.50		---
Iron	76.2	E	J	30.8	E	J	579.0	E	J	41.8	E	J	51.9	E	J	23.4	E	J	67.5	E	J
Lead	0.37		---	0.10	B	J	0.11	B	J	0.14	W	J	0.39		---	0.05	U	---	0.28		---
Magnesium	1480.0		---	57.3	B	J	162.0	B	J	251.0		---	1530.0		---	250.0		---	375.0		---
Manganese	0.38	B	J	0.23	B	J	2.70		---	0.21	B	J	0.70		---	0.25	B	J	27.30		---
Mercury	0.08		U	0.07	B	U	0.64		---	0.16		U	0.06	B	U	0.12		U	0.06	B	U
Nickel	8.20		---	0.90	U	---	0.86	U	---	0.89	U	---	0.82	U	---	0.89	U	---	0.80	U	---
Potassium	1340		---	352		---	1780		---	3450		---	1780		---	3210		---	2470		---
Selenium	0.42	U	---	0.40	U	---	0.41	U	---	0.37	U	---	0.46	U	---	0.47	U	---	0.43	U	---
Silver	0.16	U	---	0.16	U	---	0.15	U	---	0.16	U	---	0.14	U	---	0.16	U	---	0.14	U	---
Sodium	3080.0		---	759.0		---	1500.0		---	542.0		---	2650.0		---	680.0		---	1150.0		---
Thallium	0.04	UW	UJ	0.04	UW	UJ	0.04	UW	UJ	0.04	UW	UJ	0.05	UW	UJ	0.23	UW	UJ	0.04	UW	UJ
Vanadium	0.17	U	---	0.16	U	---	0.28	B	---	0.16	U	---	0.15	U	---	0.16	U	---	0.14	U	---
Zinc	80.0		---	5.1		---	25.1		---	39.4		---	57.2		---	32.5		---	23.4		---

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	Table A-7.1 Metals																				
	4K-SM-C-I	Q*	DV**	4L-SM-C-I	Q*	DV**	6A-SM-B-I	Q*	DV**	6A-SM-M-I	Q*	DV**	6B-SM-B-I	Q*	DV**	6B-SM-M-I	Q*	DV**	6C-SM-B-I	Q*	DV**
Aluminum	17.90		---	4.00	B	J	8.00	B	J	2.10	B	J	3.10	B	J	2.40	B	J	3.00	B	J
Antimony	2.7	U	---	2.3	U	---	2.6	U	---	2.7	U	---	2.5	U	---	2.6	U	---	2.7	U	---
Arsenic	0.45	U	---	0.37	U	---	0.49	UW	UJ	0.47	U	---	0.75	BW	J	0.45	U	---	0.49	UW	UJ
Barium	9.40	B	J	0.97	B	J	262.00		---	0.29	B	U	77.40		---	0.34	B	J	59.40		---
Beryllium	0.05	B	U	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---
Cadmium	0.13	U	---	0.11	U	---	0.13	U	---	0.13	U	---	0.12	U	---	0.13	U	---	0.13	U	---
Calcium	8290		---	2860		---	61600		---	184	B	J	29000		---	199	B	J	22500		---
Chromium	0.40	B	U	0.30	B	U	1.20		---	0.28	B	J	0.65		---	0.20	B	J	0.33	B	J
Cobalt	0.29	U	---	0.25	U	---	0.28	U	---	0.29	U	---	0.27	U	---	0.29	U	---	0.29	U	---
Copper	5.20		---	1.50		---	1.50		---	1.90		---	1.50		---	1.70		---	1.20	B	J
Iron	61.8	E	J	43.1	E	J	222.0		---	20.4		---	32.5		---	15.7		---	30.8		---
Lead	0.36		---	0.07	BW	J	0.69	S	---	0.08	BW	J	391.00		---	0.15	W	J	0.33		---
Magnesium	355.0		---	218.0		---	1260.0		---	302.0		---	678.0		---	318.0		---	673.0		---
Manganese	25.40		---	2.50		---	2.60		---	0.26	B	J	1.00		---	0.31	B	J	4.40		---
Mercury	0.07	B	U	0.04	B	U	0.04	U	---	0.05	U	---	0.04	U	---	0.05	U	---	0.05	U	---
Nickel	0.98	U	---	0.82	U	---	0.93	U	---	0.97	U	---	0.91	U	---	0.95	U	---	0.98	U	---
Potassium	3060		---	2260		---	1990		---	3700		---	2860		---	3780		---	2870		---
Selenium	0.45	U	---	0.37	U	---	0.05	UW	UJ	0.47	U	---	0.48	U	---	0.45	U	---	0.49	U	---
Silver	0.17	U	---	0.14	U	---	0.16	U	---	0.17	U	---	0.16	U	---	0.17	U	---	0.17	U	---
Sodium	1270.0		---	869.0		---	1870.0		---	465.0		---	1110.0		---	435.0		---	1060.0		---
Thallium	0.04	UW	UJ	0.04	UW	UJ	0.05	UW	UJ	0.05	UW	UJ	0.05	UW	UJ	0.04	UW	UJ	0.05	UW	UJ
Vanadium	0.18	U	---	0.15	U	---	0.17	U	---	0.18	U	---	0.16	U	---	0.17	U	---	0.18	U	---
Zinc	25.7		---	18.8		---	53.0	E	J	13.1	E	J	57.1	E	J	67.0	E	J	38.9	E	J

Compound (mg/Kg)	Table A-7.1 Metals																							
	6C-SM-M-I	Q*	DV**	6D-SM-B-I	Q*	DV**	6D-SM-M-I	Q*	DV**	6E-SM-B-I	Q*	DV**	6E-SM-M-I	Q*	DV**	6F-SM-B-I	Q*	DV**	6F-SM-M-I	Q*	DV**
Aluminum	1.80	B	J	3.50	B	J	1.30	B	J	3.70	B	J	0.97	U	---	3.30	B	J	1.60	B	J
Antimony	2.5	U	---	2.6	U	---	2.6	U	---	2.6	U	---	2.6	U	---	2.6	U	---	2.6	U	---
Arsenic	0.48	U	---	0.48	UW	UJ	0.49	U	---	0.49	U	---	0.45	U	---	0.47	UW	UJ	0.42	U	---
Barium	0.34	B	---	168.00		---	0.33	B	U	113.00		---	0.30	B	U	53.70		---	0.30	B	U
Beryllium	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.04	U	---
Cadmium	0.12	U	---	0.13	U	---	0.13	U	---	0.12	U	---	0.13	U	---	0.12	U	---	0.13	U	---
Calcium	174	B	J	56700		---	191	B	J	36700		---	166	B	J	44200		---	193	B	---
Chromium	0.22	B	J	0.84		---	0.25	B	J	0.82		---	0.23	B	J	0.95		---	0.19	B	---
Cobalt	0.27	U	---	0.28	U	---	0.28	U	---	0.28	U	---	0.28	U	---	0.28	U	---	0.28	U	---
Copper	1.60		---	1.70		---	1.10	B	J	6.20		---	1.40		---	9.60		---	1.40		---
Iron	14.2		---	117.0		---	16.5		---	35.2		---	15.8		---	37.4		---	21.0		---
Lead	0.75		---	5.60		---	0.37		---	49.70		---	0.14		---	0.26	W	J	0.23	W	J
Magnesium	285.0		---	1040.0		---	310.0		---	818.0		---	299.0		---	933.0		---	319.0		---
Manganese	2.00		---	2.00		---	0.39	B	J	1.60		---	0.29	B	J	2.30		---	0.49	B	J
Mercury	0.04	B	U	0.03	B	U	0.04	U	---	0.04	U	---	0.05	B	U	0.04	U	---	0.04	U	---
Nickel	0.89	U	---	0.93	U	---	0.93	U	---	0.92	U	---	0.93	U	---	0.92	U	---	0.94	U	---
Potassium	3490		---	1580		---	3830		---	2870		---	3620		---	2760		---	3980		---
Selenium	0.48	U	---	0.07	BW	J	0.49	U	---	0.49	U	---	0.45	U	---	0.47	U	---	0.42	U	---
Silver	0.16	U	---	0.16	U	---	0.16	U	---	0.16	U	---	0.16	U	---	0.16	U	---	0.16	U	---
Sodium	482.0		---	1300.0		---	421.0		---	1200.0		---	363.0		---	1420.0		---	428.0		---
Thallium	0.05	UW	UJ	0.05	U	---	0.05	UW	UJ	0.05	UW	UJ	0.04	UW	UJ	0.05	UW	UJ	0.04	UW	UJ
Vanadium	0.16	U	---	0.17	U	---	0.17	U	---	0.17	U	---	0.17	U	---	0.17	U	---	0.17	U	---
Zinc	22.8	E	J	44.8	E	J	11.3	E	J	94.0	E	J	11.1	E	J	46.8	E	J	20.1	E	J

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	Table A-7.1 Metals					
	6G-SM-B-I	Q*	DV**	6G-SM-M-I	Q*	DV**
Aluminum	2.80	B	J	1.50	B	J
Antimony	2.5	U	---	2.7	U	---
Arsenic	0.45	UW	UJ	0.45	U	---
Barium	225.00		---	0.47	B	
Beryllium	0.04	U	---	0.04	U	---
Cadmium	0.12	U	---	0.13	U	---
Calcium	66200		---	226	B	J
Chromium	5.10		---	0.18	B	J
Cobalt	0.27	U	---	0.29	U	---
Copper	0.78	B	U	1.30		---
Iron	56.4	E	J	55.2		---
Lead	43.70		---	19.30		---
Magnesium	1300.0		---	288.0		---
Manganese	4.60		---	0.22	B	U
Mercury	0.05	B	U	0.04	U	---
Nickel	2.40		---	0.96	U	---
Potassium	2210		---	3700		---
Selenium	0.04	UW	UJ	0.45	U	---
Silver	0.16	U	---	0.17	U	---
Sodium	1550.0		---	577.0		---
Thallium	0.04	UW	UJ	0.05	UW	UJ
Vanadium	0.16	U	---	0.17	U	---
Zinc	51.1		---	9.3	E	UJ

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-7.2 Pesticides/PCB's																							
	2A-SM-B-I	Q*	DV**	2A-SM-F-I	Q*	DV**	2A-SM-L-I	Q*	DV**	2A-SM-M-I	Q*	DV**	3A-SM-B-I	Q*	DV**	3A-SM-M-I	Q*	DV**	4H-SM-C-I	Q*	DV**
Alpha-BHC	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Beta-BHC	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Delta-BHC	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Gamma-BHC	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Heptachlor	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Aldrin	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Heptachlor epoxide	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Endosulfan I	47	U	---	48	U	---	48	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Dieldrin	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
4,4'-DDE	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
Endrin	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
Endosulfan II	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
4,4'-DDD	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
Endosulfan sulfate	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
4,4'-DDT	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
Methoxychlor	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Endrin ketone	94	U	---	95	U	---	96	U	---	94	U	---	95	U	---	95	U	---	94	U	---
Alpha-chlordane	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Gamma-chlordane	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Toxaphene	940	U	---	950	U	---	960	U	---	940	U	---	950	U	---	950	U	---	940	U	---
Aroclor-1016	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1221	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-7.2 Pesticides/PCB's																							
	2A-SM-B-I	Q*	DV**	2A-SM-F-I	Q*	DV**	2A-SM-L-I	Q*	DV**	2A-SM-M-I	Q*	DV**	3A-SM-B-I	Q*	DV**	3A-SM-M-I	Q*	DV**	4H-SM-C-I	Q*	DV**
Aroclor-1232	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1242	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1248	470	U	---	480	U	---	480	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1254	940	U	---	950	U	---	960	U	---	940	U	---	950	U	---	950	U	---	940	U	---
Aroclor-1260	940	U	---	950	U	---	960	U	---	940	U	---	950	U	---	950	U	---	940	U	---

Compound (ug/Kg)	Table A-7.2 Pesticides/PCB's																							
	4K-SM-C-I	Q*	DV**	4L-SM-C-I	Q*	DV**	6A-SM-B-I	Q*	DV**	6A-SM-M-I	Q*	DV**	6B-SM-B-I	Q*	DV**	6B-SM-M-I	Q*	DV**	6C-SM-B-I	Q*	DV**
Alpha-BHC	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Beta-BHC	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Delta-BHC	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Gamma-BHC	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Heptachlor	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Aldrin	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Heptachlor epoxide	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Endosulfan I	47	U	---	46	U	---	47	U	---	47	U	---	47	U	---	47	U	---	46	U	---
Dieldrin	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
4,4'-DDE	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
Endrin	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
Endosulfan II	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
4,4'-DDD	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
Endosulfan sulfate	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
4,4'-DDT	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
Methoxychlor	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---
Endrin ketone	94	U	---	92	U	---	95	U	---	94	U	---	93	U	---	95	U	---	93	U	---
Alpha-chlordane	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---
Gamma-chlordane	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---
Toxaphene	940	U	---	920	U	---	950	U	---	940	U	---	930	U	---	950	U	---	930	U	---
Aroclor-1016	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---
Aroclor-1221	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---

Compound (ug/Kg)	Table A-7.2 Pesticides/PCB's																							
	4K-SM-C-I	Q*	DV**	4L-SM-C-I	Q*	DV**	6A-SM-B-I	Q*	DV**	6A-SM-M-I	Q*	DV**	6B-SM-B-I	Q*	DV**	6B-SM-M-I	Q*	DV**	6C-SM-B-I	Q*	DV**
Aroclor-1232	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---
Aroclor-1242	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---
Aroclor-1248	470	U	---	460	U	---	470	U	---	470	U	---	470	U	---	470	U	---	460	U	---
Aroclor-1254	940	U	---	920	U	---	950	U	---	940	U	---	930	U	---	950	U	---	930	U	---
Aroclor-1260	940	U	---	920	U	---	950	U	---	940	U	---	930	U	---	950	U	---	930	U	---

Compound (ug/Kg)	Table A-7.2 Pesticides/PCB's																							
	6C-SM-M-I	Q*	DV**	6D-SM-B-I	Q*	DV**	6D-SM-M-I	Q*	DV**	6E-SM-B-I	Q*	DV**	6E-SM-M-I	Q*	DV**	6F-SM-B-I	Q*	DV**	6F-SM-M-I	Q*	DV**
Alpha-BHC	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Beta-BHC	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Delta-BHC	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Gamma-BHC	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Heptachlor	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Aldrin	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Heptachlor epoxide	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Endosulfan I	47	U	---	46	U	---	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Dieldrin	93	U	---	93	U	---	94	U	---	93	U	---	96	U	---	93	U	---	94	U	---
4,4'-DDE	93	U	---	93	U	---	94	U	---	93	U	---	96	U	---	93	U	---	94	U	---
Endrin	93	U	---	93	U	---	94	U	---	93	U	---	180	U		93	U	---	94	U	---
Endosulfan II	93	U	---	93	U	---	94	U	---	93	U	---	96	U	---	93	U	---	94	U	---
4,4'-DDD	93	U	---	93	U	---	94	U	---	93	U	---	110	U	J	93	U	---	94	U	---
Endosulfan sulfate	93	U	---	93	U	---	94	U	---	93	U	---	96	U	---	93	U	---	94	U	---
4,4'-DDT	93	U	---	93	U	---	94	U	---	93	U	---	96	U	---	93	U	---	94	U	---
Methoxychlor	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Endrin ketone	93	U	---	93	U	---	94	U	---	93	U	---	96	U	---	93	U	---	94	U	---
Alpha-chlordane	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Gamma-chlordane	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Toxaphene	930	U	---	930	U	---	940	U	---	930	U	---	960	U	---	930	U	---	940	U	---
Aroclor-1016	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1221	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---

Compound (ug/Kg)	Table A-7.2 Pesticides/PCB's																							
	6C-SM-M-I	Q*	DV**	6D-SM-B-I	Q*	DV**	6D-SM-M-I	Q*	DV**	6E-SM-B-I	Q*	DV**	6E-SM-M-I	Q*	DV**	6F-SM-B-I	Q*	DV**	6F-SM-M-I	Q*	DV**
Aroclor-1232	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1242	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1248	470	U	---	460	U	---	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1254	930	U	---	930	U	---	940	U	---	930	U	---	960	U	---	930	U	---	940	U	---
Aroclor-1260	930	U	---	930	U	---	940	U	---	930	U	---	960	U	---	930	U	---	940	U	---

	Table A-7.2 Pesticides/PCB's					
Compound (ug/Kg)	6G-SM-B-I	Q*	DV**	6G-SM-M-I	Q*	DV**
Alpha-BHC	48	U	---	47	U	---
Beta-BHC	48	U	---	47	U	---
Delta-BHC	48	U	---	47	U	---
Gamma-BHC	48	U	---	47	U	---
Heptachlor	48	U	---	47	U	---
Aldrin	48	U	---	47	U	---
Heptachlor epoxide	48	U	---	47	U	---
Endosulfan I	48	U	---	47	U	---
Dieldrin	95	U	---	94	U	---
4,4'-DDE	95	U	---	94	U	---
Endrin	95	U	---	94	U	---
Endosulfan II	95	U	---	94	U	---
4,4'-DDD	95	U	---	94	U	---
Endosulfan sulfate	95	U	---	94	U	---
4,4'-DDT	95	U	---	94	U	---
Methoxychlor	480	U	---	470	U	---
Endrin ketone	95	U	---	94	U	---
Alpha-chlordane	480	U	---	470	U	---
Gamma-chlordane	480	U	---	470	U	---
Toxaphene	950	U	---	940	U	---
Aroclor-1016	480	U	---	470	U	---
Aroclor-1221	480	U	---	470	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-7.2 Pesticides/PCB's					
	6G-SM-B-I	Q*	DV**	6G-SM-M-I	Q*	DV**
Aroclor-1232	480	U	---	470	U	---
Aroclor-1242	480	U	---	470	U	---
Aroclor-1248	480	U	---	470	U	---
Aroclor-1254	950	U	---	940	U	---
Aroclor-1260	950	U	---	940	U	---

Parameter (pCi/g)	Table A-7.3 Gross Alpha/Beta																			
	2A-SM-B-I	DV**	2A-SM-F-I	DV**	2A-SM-L-I	DV**	2A-SM-M-I	DV**	3A-SM-B-I	DV**	3A-SM-M-I	DV**	4H-SM-C-I	DV**	4K-SM-C-I	DV**	4L-LM-C-I	DV**
Alpha	5.24	---	ND	---	ND	---	0.09	---	ND	---	ND	---	ND	---	0.09	---	0.09	---
Sigma Error +/-	6.84	---	-	---	-	---	0.11	---	-	---	-	---	-	---	0.08	---	0.08	---
MDA*	11.46	U	0.15	U	0.36	U	0.17	U	13.06	U	0.15	U	0.21	U	0.11	U	0.12	U
Beta	18.90	---	0.09	---	1.57	---	3.14	---	19.43	---	3.14	---	0.82	---	0.89	---	0.61	---
Sigma Error +/-	5.42	---	0.08	---	0.15	---	0.18	---	5.25	---	0.16	---	0.11	---	0.10	---	0.11	---
MDA*	7.56	---	0.13	U	0.13	---	0.12	---	7.44	---	0.10	---	0.13	---	0.11	---	0.11	---

Parameter (pCi/g)	Table A-7.3 Gross Alpha/Beta																			
	6A-SM-B-I	DV**	6A-SM-M-I	DV**	6B-SM-B-I	DV**	6B-SM-M-I	DV**	6C-SM-B-I	DV**	6C-SM-M-I	DV**	6D-SM-B-I	DV**	6D-SM-M-I	DV**	6E-SM-B-I	DV**
Alpha	5.96	---	ND	---	12.55	J	1.96	---	ND	---	6.44	---	4.32	---	0.19	---	ND	---
Sigma Error +/-	5.35	---	-	---	7.86	---	0.37	---	-	---	0.58	---	5.61	---	0.13	---	-	---
MDA*	7.97	UJ	0.31	U	10.06	---	0.24	---	11.38	UJ	0.16	---	9.30	UJ	0.19	---	9.84	U
Beta	52.10	---	2.92	U	107.27	---	3.28	U	85.70	---	3.11	U	57.22	---	2.83	U	71.91	---
Sigma Error +/-	5.78	---	0.19	---	8.78	---	0.10	---	7.77	---	0.18	---	7.11	---	0.17	---	7.08	---
MDA*	5.70	---	0.13	---	7.03	---	0.12	---	7.35	---	0.12	---	7.47	---	0.13	---	6.73	---

Parameter (pCi/g)	Table A-7.3 Gross Alpha/Beta									
	6E-SM-M-I	DV**	6F-SM-B-I	DV**	6F-SM-M-I	DV**	6G-SM-B-I	DV**	6G-SM-M-I	DV**
Alpha	ND	---	ND	---	0.20	---	ND	---	0.14	---
Sigma Error +/-	-	---	-	---	0.14	---	-	---	0.13	---
MDA*	0.28	U	0.25	UJ	0.21	U	14.13	U	0.20	UJ
Beta	2.65	U	1.03	U	2.60	U	59.80	---	1.80	---
Sigma Error +/-	0.17	---	0.12	---	0.16	---	7.74	---	0.15	---
MDA*	0.14	---	0.12	---	0.10	---	8.53	---	0.13	---

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy																	
	2A-SM-B-I	DV**	2A-SM-F-I	DV**	2A-SM-L-I	DV**	2A-SM-M-I	DV**	3A-SM-B-I	DV**	3A-SM-L-I	DV**	3A-SM-M-I	DV**	4A-SM-C-I	DV**	4B-SM-C-I	DV**
Potassium 40	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	2.90	U	2.20	U	4.20	U	3.00	U	4.30	U	5.30	U	3.10	U	4.20	U	2.90	U
Cesium 137	0.14	---	ND	---	0.59	---	ND	---	0.28	---	0.60	---	0.63	---	0.53	---	0.27	---
Sigma Error +/-	0.03	---	-	---	0.06	---	-	---	0.04	---	0.10	---	0.05	---	0.06	---	0.04	---
MDA*	0.10	---	0.10	U	0.10	---	0.20	U	0.10	---	0.10	---	0.10	---	0.10	---	0.10	---
Lead 212	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.10	U	0.10	U	0.30	U	0.10	U	0.20	U	0.20	U	0.20	U	0.30	U	0.20	U
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.30	U	0.60	U	0.20	U	0.30	U	0.30	U	0.20	U	0.40	U	0.40	U
Uranium 235	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.50	U	0.40	U	1.00	U	0.50	U	0.60	U	0.60	U	0.40	U	0.80	U	0.70	U
Uranium238	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.40	U	0.90	U	1.90	U	1.30	U	1.20	U	1.30	U	1.40	U	1.90	U	1.60	U
Cerium 144	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	U	0.30	U	0.80	U	0.50	U	0.40	U	0.50	U	0.40	U	0.80	U	0.50	U
Thallium 208	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.10	U	0.10	U	0.30	U	0.10	U	0.10	U	0.10	U	0.10	U	0.20	U	0.20	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy																	
	2A-SM-B-I	DV**	2A-SM-F-I	DV**	2A-SM-L-I	DV**	2A-SM-M-I	DV**	3A-SM-B-I	DV**	3A-SM-L-I	DV**	3A-SM-M-I	DV**	4A-SM-C-I	DV**	4B-SM-C-I	DV**
Radium 226	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.10	U	0.50	U	0.20	U	0.30	U	0.30	U	0.20	U	0.40	U	0.30	U
Cerium 139	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.06	U	0.04	U	0.10	U	0.06	U	0.07	U	0.08	U	0.06	U	0.10	U	0.08	U
Lead 210	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	10.70	U	0.05	U	1.90	U	10.40	U	1.40	U	1.60	U	14.50	U	1.90	U	0.08	U
Thorium 230	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

Table A-7.4 Gamma Spectroscopy																				
Nuclide (pCi/g)	4C-SM-C-I	DV**	4D-SM-C-I	DV**	4E-SM-C-I	DV**	4F-SM-C-I	DV**	4G-SM-C-I	DV**	4H-SM-C-I	DV**	4I-SM-C-I	DV**	4J-SM-C-I	DV**	4K-SM-C-I	DV**
Potassium 40	ND	---	2.80	---	4.40	---	ND	---	4.90	---	ND	---	ND	---	ND	---	3.60	---
Sigma Error +/-	-	---	0.60	---	1.20	---	-	---	0.90	---	-	---	-	---	-	---	0.40	---
MDA*	3.70	U	2.40	---	4.20	---	2.60	U	3.50	---	0.20	U	3.20	U	6.60	U	1.90	---
Cesium 137	0.18	---	ND	---	ND	---	ND	---	0.33	---	ND	---	0.26	---	ND	---	3.50	---
Sigma Error +/-	0.03	---	-	---	-	---	-	---	0.04	---	-	---	0.03	---	-	---	0.20	---
MDA*	0.10	---	0.10	U	0.20	U	0.10	U	0.10	---	0.20	U	0.10	---	0.10	U	0.30	---
Lead 212	ND	---	ND	---	0.40	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	0.10	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	0.20	U	0.30	---	0.20	U	0.20	U	0.10	U	0.20	U	0.30	U	0.20	U
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	U	0.30	U	0.40	U	0.20	U	0.30	U	0.30	U	0.30	U	0.30	U	0.30	U
Uranium 235	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.70	U	0.40	U	0.80	U	0.70	U	0.70	U	0.40	U	0.60	U	1.00	U	0.50	U
Uranium238	ND	---	ND	---	1.90	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	0.05	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.50	U	1.30	U	1.70	---	1.80	U	1.40	U	1.30	U	1.50	U	1.80	U	1.10	U
Cerium 144	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.60	U	0.40	U	0.70	U	0.50	U	0.60	U	0.40	U	0.60	U	0.80	U	0.30	U
Thallium 208	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.10	U	0.20	U	0.10	U	0.20	U	0.10	U	0.20	U	0.20	U	0.10	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy																	
	4C-SM-C-I	DV**	4D-SM-C-I	DV**	4E-SM-C-I	DV**	4F-SM-C-I	DV**	4G-SM-C-I	DV**	4H-SM-C-I	DV**	4I-SM-C-I	DV**	4J-SM-C-I	DV**	4K-SM-C-I	DV**
Radium 226	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	0.20	U	0.30	U	0.20	U	0.30	U	0.20	U	0.30	U	0.30	U	0.20	U
Cerium 139	ND	---	ND	---	0.14	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	0.02	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.10	U	0.07	U	0.10	---	0.09	U	0.10	U	0.60	U	0.09	U	0.10	U	0.06	U
Lead 210	ND	---	ND	---	1.70	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	0.40	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.90	U	0.03	U	1.50	---	15.70	U	1.70	U	10.50	U	1.50	U	2.00	U	0.07	U
Thorium 230	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy																	
	4L-SM-C-I	DV**	4M-SM-C-I	DV**	6A-SM-B-I	DV**	6A-SM-L-I	DV**	6A-SM-M-I	DV**	6B-SM-B-I	DV**	6B-SM-L-I	DV**	6B-SM-M-I	DV**	6C-SM-B-I	DV**
Potassium 40	ND	---	5.60	---	ND	---	9.50	---	ND	---	ND	---	ND	---	3.00	---	ND	---
Sigma Error +/-	-	---	1.10	---	-	---	1.40	---	-	---	-	---	-	---	0.30	---	-	---
MDA*	3.90	U	3.50	---	2.40	U	7.50	---	1.80	U	6.90	U	3.00	U	1.60	---	1.90	U
Cesium 137	ND	---	ND	---	ND	---	0.31	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	0.05	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.10	U	0.20	U	0.10	U	0.10	---	0.20	U	0.20	U	0.10	U	0.10	U	0.20	U
Lead 212	ND	---	ND	---	0.54	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	0.05	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.20	U	0.20	---	0.30	U	0.10	U	0.30	U	0.20	U	0.20	U	0.10	U
Bismuth 214	ND	---	-	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	-	---	0.20	U	0.40	U	0.30	U	0.50	U	0.30	U	0.20	U	0.20	U
Uranium 235	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	U	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Uranium238	ND	---	-	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.10	U	-	---	1.10	U	2.00	U	1.40	U	1.80	U	1.20	U	1.30	U	1.10	U
Cerium 144	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Thallium 208	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	0.13	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	0.03	---	-	---	-	---
MDA*	0.10	U	0.20	U	0.10	U	0.20	U	0.10	U	0.20	U	0.10	---	0.10	U	0.10	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy																	
	4L-SM-C-I	DV**	4M-SM-C-I	DV**	6A-SM-B-I	DV**	6A-SM-L-I	DV**	6A-SM-M-I	DV**	6B-SM-B-I	DV**	6B-SM-L-I	DV**	6B-SM-M-I	DV**	6C-SM-B-I	DV**
Radium 226	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.30	U	0.20	U	0.40	U	0.20	U	0.40	U	0.30	U	0.20	U	0.20	U
Cerium 139	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.06	U	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Lead 210	ND	---	-	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.07	U	-	---	1.10	U	2.40	U	10.10	U	2.30	U	1.20	U	10.00	U	0.10	U
Thorium 230	-	---	-	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	14.50	U	23.80	U	38.30	U	21.60	U	14.90	U	36.20	U	11.40	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	6C-SM-L-I	DV**	6C-SM-M-I	DV**	6D-SM-B-I	DV**	6D-SM-L-I	DV**	6D-SM-M-I	DV**	6E-SM-B-I	DV**	6E-SM-L-I	DV**	6E-SM-M-I	DV**	6F-SM-B-I	DV**
Table A-7.4 Gamma Spectroscopy																		
Potassium 40	ND	---	ND	---	ND	---	ND	---	4.20	---	2.70	---	5.20	---	2.50	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	0.50	---	0.30	---	0.80	---	0.40	---	-	---
MDA*	5.60	U	1.60	U	2.50	U	3.30	U	1.90	---	2.60	---	3.20	---	1.80	---	6.60	U
Cesium 137	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	0.10	U	0.10	U	0.20	U	0.10	U	0.10	U	0.20	U	0.10	U	0.20	U
Lead 212	0.80	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	0.10	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	---	0.10	U	0.20	U	0.20	U	0.20	U	0.20	U	0.20	U	0.10	U	0.30	U
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.60	U	0.20	U	0.20	U	0.40	U	0.30	U	0.20	U	0.30	U	0.20	U	0.40	U
Uranium 235	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Uranium238	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	2.60	U	1.10	U	1.10	U	1.40	U	1.10	U	1.10	U	1.30	U	1.40	U	1.80	U
Cerium 144	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Thallium 208	ND	---	ND	---	ND	---	0.14	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	0.04	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.10	U	0.20	U	0.10	---	0.10	U	0.10	U	0.10	U	0.10	U	0.20	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy																	
	6C-SM-L-I	DV**	6C-SM-M-I	DV**	6D-SM-B-I	DV**	6D-SM-L-I	DV**	6D-SM-M-I	DV**	6E-SM-B-I	DV**	6E-SM-L-I	DV**	6E-SM-M-I	DV**	6F-SM-B-I	DV**
Radium 226	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.50	U	0.20	U	0.20	U	0.30	U	0.30	U	0.20	U	0.30	U	0.20	U	0.40	U
Cerium 139	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Lead 210	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	2.50	U	9.10	U	1.40	U	1.80	U	0.10	U	1.00	U	1.60	U	9.60	U	1.60	U
Thorium 230	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	30.60	U	33.60	U	14.30	U	16.60	U	10.70	U	12.90	U	15.00	U	33.10	U	21.10	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy									
	6F-SM-L-I	DV**	6F-SM-M-I	DV**	6G-SM-B-I	DV**	6G-SM-L-I	DV**	6G-SM-M-I	DV**
Potassium 40	9.80	---	3.10	---	ND	---	1.90	---	3.90	---
Sigma Error +/-	1.20	---	0.40	---	-	---	0.40	---	0.40	---
MDA*	7.70	---	1.80	---	6.50	U	1.68	---	1.70	---
Cesium 137	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.10	U	0.20	U	0.10	U	0.10	U
Lead 212	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	U	0.20	U	0.30	U	0.10	U	0.20	U
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	U	0.20	U	0.50	U	0.20	U	0.20	U
Uranium 235	-	---	-	---	ND	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	0.10	U	-	---	-	---
Uranium238	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	2.20	U	1.10	U	1.80	U	1.00	U	1.40	U
Cerium 144	-	---	-	---	ND	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	0.90	U	-	---	-	---
Thallium 208	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.10	U	0.20	U	0.10	U	0.10	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-7.4 Gamma Spectroscopy									
	6F-SM-L-I	DV**	6F-SM-M-I	DV**	6G-SM-B-I	DV**	6G-SM-L-I	DV**	6G-SM-M-I	DV**
Radium 226	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	U	0.20	U	0.40	U	0.20	U	0.20	U
Cerium 139	-	---	-	---	ND	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	0.10	U	-	---	-	---
Lead 210	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---
MDA*	2.00	U	10.40	U	2.20	U	0.03	U	9.70	U
Thorium 230	22.80	---	ND	---	-	---	ND	---	ND	---
Sigma Error +/-	4.40	---	-	---	-	---	-	---	-	---
MDA*	18.90	---	36.00	U	-	---	9.30	U	36.00	U

* Minimum Detectable Activity. ** Data Validation.

Table A-7.5 Tritium

Nuclide (pCi/g)	2A-SM-B-I	DV**	2A-SM-F-I	DV**	2A-SM-L-I	DV**	2A-SM-M-I	DV**	3A-SM-B-I	DV**	3A-SM-L-I	DV**	3A-SM-M-I	DV**	4A-SM-C-I	DV**	4B-SM-C-I	DV**
Tritium	12.70	---	0.10	---	ND	---	ND	---	0.35	---	1.14	---	6.37	---	ND	---	0.83	---
Sigma Error +/-	1.26	---	0.03	---	-	---	-	---	0.06	---	0.15	---	0.65	---	-	---	0.10	---
MDA*	0.07	---	0.07	---	0.13	U	0.13	U	0.08	---	0.14	---	0.12	---	0.14	U	0.09	---

Nuclide (pCi/g)	4C-SM-C-I	DV**	4D-SM-C-I	DV**	4E-SM-C-I	DV**	4F-SM-C-I	DV**	4G-SM-C-I	DV**	4H-SM-C-I	DV**	4I-SM-C-I	DV**	4J-SM-C-I	DV**	4K-SM-C-I	DV**
Table A-7.5 Tritium																		
Tritium	3.85	---	0.45	---	39.77	---	0.20	---	1.46	---	0.18	---	0.08	---	0.59	---	0.51	---
Sigma Error +/-	0.41	---	0.08	---	3.93	---	0.07	---	0.17	---	0.07	---	0.05	---	0.10	---	0.09	---
MDA*	0.14	---	0.13	---	0.16	---	0.14	---	0.12	---	0.14	---	0.12	U	0.14	---	0.14	---

* Minimum Detectable Activity. **Data Validation.

Table A-7.5 Tritium

Nuclide (pCi/g)	4L-SM-C-I	DV**	4M-SM-C-I	DV**	6A-SM-B-I	DV**	6A-SM-L-I	DV**	6A-SM-M-I	DV**	6B-SM-B-I	DV**	6B-SM-L-I	DV**	6B-SM-M-I	DV**	6C-SM-B-I	DV**
Tritium	0.39	---	0.25	---	ND	---	ND	---	ND	---	ND	---	20.17	---	11.38	---	ND	---
Sigma Error +/-	0.08	---	0.07	---	-	---	-	---	-	---	-	---	2.00	---	1.14	---	-	---
MDA*	0.14	---	0.14	---	0.18	U	0.22	U	0.23	U	0.15	U	0.22	---	0.21	---	0.12	U

Table A-7.5 Tritium																						
Nuclide (pCi/g)	6C-SM-L-I	DV**	6C-SM-M-I	DV**	6D-SM-B-I	DV**	6D-SM-L-I	DV**	6D-SM-M-I	DV**	6E-SM-B-I	DV**	6E-SM-L-I	DV**	6E-SM-M-I	DV**	6F-SM-B-I	DV**
Tritium	ND	---	ND	---	0.25	---	ND	---	ND	---	0.38	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	0.07	---	-	---	-	---	0.06	---	-	---	-	---	-	---
MDA*	0.29	U	0.20	U	0.13	---	0.16	U	0.19	U	0.09	---	0.11	---	0.05	---	0.10	U

Nuclide (pCi/g)	Table A-7.5 Tritium									
	6F-SM-L-I	DV**	6F-SM-M-I	DV**	6G-SM-B-I	DV**	6G-SM-M-I	DV**	6G-SM-L-I	DV**
Tritium	ND	---	ND	---	ND	---	ND	---	0.12	---
Sigma Error +/-	-	---	-	---	-	---	-	---	0.06	---
MDA*	0.13	---	0.13	U	0.14	U	0.12	---	0.13	U

Compound (mg/Kg)	Table A-8.1 Metals																					
	1A-LM-B-I	Q*	DV**	1A-LM-L-I	Q*	DV**	1A-LM-M-I	Q*	DV**	1B-LM-B-I	Q*	DV**	1B-LM-F-I	Q*	DV**	1B-LM-L-I	Q*	DV**	1B-LM-M-I	Q*	DV**
Aluminum	7.4	B	J	2.5	B	U	1.5	B	U	5.8	B	U	2.1	B	U	1.8	B	U	1.9	B	U
Antimony	2.6	U	---	4.7	U	---	2.4	U	---	2.6	U	---	2.7	U	---	4.5	U	---	2.5	U	---
Arsenic	0.48	UW	UJ	0.07	U	---	0.05	UW	UJ	0.46	UW	UJ	0.05	UW	UJ	0.07	U	---	0.04	UW	UJ
Barium	370.00		---	0.94	B	J	0.39	B	J	345.00		---	0.80	B	J	0.61	B	J	0.28	B	J
Beryllium	0.04	U	---	0.08	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.07	U	---	0.04	U	---
Cadmium	0.13	U	---	0.23	U	---	0.12	U	---	0.15	B	J	0.13	U	---	0.22	U	---	0.12	U	---
Calcium	209000		---	440		---	163	B	J	199000		---	380		---	355	B	J	185	B	J
Chromium	3.40		---	0.58	B	J	0.22	B	J	3.20		---	0.61		---	0.39	B	J	0.47		---
Cobalt	0.28	U	---	0.50	U	---	0.26	U	---	0.28	U	---	0.29	U	---	0.48	U	---	0.27	U	---
Copper	0.28	B	U	62.60		---	1.50		---	0.17	U	---	3.00		---	33.50		---	1.70		---
Iron	10.4		---	146.0		---	24.0		---	2.8	B	J	35.8		---	213.0		---	29.9		---
Lead	13.90		---	0.05	B	J	0.04	U	---	0.38		---	0.27		---	0.05	UW	UJ	0.06	U	UJ
Magnesium	3350.0		---	184.0	B	J	256.0		---	3100.0		---	71.7	B	J	199.0	B	J	270.0		---
Manganese	1.20		---	6.30		---	4.20		---	1.30		---	2.60		---	5.20		---	3.30		---
Mercury	0.04	U	---	0.04	B	J	0.07		---	0.03	U	---	0.04	U	---	0.06	B	J	0.19		---
Nickel	0.94	U	---	1.70	U	---	0.88	U	---	0.94	U	---	0.97	U	---	1.60	U	---	0.91	U	---
Potassium	356		U	2620		---	3380		---	368		U	821		---	3200		---	3410		---
Selenium	0.48	U	---	0.73	U	---	0.45	U	---	0.46	U	---	0.47	U	---	0.70	U	---	0.43	U	---
Silver	0.16	U	---	0.29	U	---	0.15	U	---	0.16	U	---	0.17	U	---	0.28	U	---	0.16	U	---
Sodium	5190.0		---	925.0		---	735.0		---	5220.0		---	527.0		---	906.0		---	635.0		U
Thallium	0.05	UW	UJ	0.07	UW	UJ	0.05	UW	UJ	0.05	U	---	0.05	UW	UJ	0.07	UW	UJ	0.04	UW	UJ
Vanadium	0.17	U	---	0.30	U	---	0.16	U	---	0.17	U	---	0.17	U	---	0.29	U	---	0.16	U	---
Zinc	99.1		---	32.2		---	31.3		---	70.4		---	7.6		---	39.8		---	43.7		---

* Data Qualifier. ** Data Validation.

Table A-8.1 Metals

Compound (mg/Kg)	1C-LM-B-I	Q*	DV**	1C-LM-B-I-D	Q*	DV**	1C-LM-F-I	Q*	DV**	1C-LM-F-I-D	Q*	DV**	1C-LM-L-I	Q*	DV**	1C-LM-L-I-D	Q*	DV**	1C-LM-M-I	Q*	DV**
Aluminum	5.1	B	U	4.5	B	---	3.1	B	U	2.1	B	---	1.9	B	U	1.6	U	---	1.9	B	U
Antimony	2.1	U	---	2.1	U	---	2.6	U	---	2.9	B	---	4.3	U	---	4.2	U	---	2.5	U	---
Arsenic	0.48	UE	UJ	0.48	U	---	0.05	U	---	0.05	U	---	0.06	U	---	0.09	U	---	0.05	UW	UJ
Barium	415.00		---	405.00		---	0.52	B	J	0.82	B	---	0.46	B	J	0.40	B	---	0.34	B	J
Beryllium	0.03	U	---	0.03	U	---	0.04	U	---	0.04	U	---	0.07	U	---	0.07	U	---	0.04	U	---
Cadmium	0.10	U	---	0.10	U	---	0.13	U	---	0.13	U	---	0.21	U	---	0.22	B	---	0.12	U	---
Calcium	226000		---	231000		---	174	B	J	388		---	185	B	J	246	B	---	220	B	J
Chromium	3.60		---	3.50		---	0.50		---	0.42	B	---	0.31	B	J	0.42	B	---	0.33	B	J
Cobalt	0.22	U	---	0.23	U	---	0.28	U	---	0.29	U	---	0.59	B	J	0.46	U	---	0.27	U	---
Copper	0.14	U	---	0.14	U	---	2.60		---	0.76	B	---	52.50		---	52.00		---	1.50		---
Iron	3.5	B	J	2.4	B	---	11.2		---	39.7		---	195.0		---	177.0		---	24.8		---
Lead	Not Run	-	---	Not Run	-	---	88.20		---	20.40		---	6.70	S	---	2.30		---	0.09	B	J
Magnesium	3190.0		---	3200.0		---	39.3	B	J	50.3	B	---	180.0	B	J	170.0	B	---	237.0		---
Manganese	1.40		---	1.40		---	3.20		---	0.50	B	---	5.10		---	4.80		---	0.27	B	J
Mercury	0.03	U	---	0.03	U	---	0.05	U	---	0.04	U	---	0.05	B	J	0.05	B	---	0.04	U	---
Nickel	0.75	U	---	0.77	U	---	0.93	U	---	0.96	U	---	1.60	U	---	1.50	U	---	0.90	U	---
Potassium	309		U	335		---	434		U	596		---	2470		---	2380		---	2890		---
Selenium	0.48	U	---	0.48	U	---	0.47	U	---	0.46	U	---	0.59	U	---	0.88	U	---	0.49	U	---
Silver	0.13	U	---	0.13	U	---	0.16	U	---	0.17	U	---	0.27	U	---	0.27	U	---	0.16	U	---
Sodium	5720.0		---	5590.0		---	218.0	B	J	411.0		---	1260.0		---	1184.0		---	568.0		U
Thallium	0.05	U	---	0.05	U	---	0.05	UW	UJ	0.05	U	---	0.06	UW	UJ	0.09	U	---	0.05	UW	UJ
Vanadium	0.13	U	---	0.14	U	---	0.17	U	---	0.17	U	---	0.28	U	---	0.28	U	---	0.16	U	---
Zinc	51.6		---	53.0		---	3.3		---	4.4		---	36.8		---	36.6		---	30.3		---

* Data Qualifier. ** Data Validation.

Table A-8.1 Metals

Compound (mg/Kg)	1C-LM-M-I-D	Q*	DV**	1D-LM-B-I	Q*	DV**	1D-LM-B-I-D	Q*	DV**	1D-LM-F-I	Q*	DV**	1D-LM-F-I-D	Q*	DV**	1D-LM-L-I	Q*	DV**	1D-LM-L-I-D	Q*	DV**
Aluminum	1.7	B	---	5.5	B	J	5.6	B	NP	4.3	B	J	4.0	U	NP	1.5	B	J	1.6	B	NP
Antimony	2.5	U	---	2.2	U	---	2.4	U	NP	9.7	U	---	10.5	U	NP	2.6	U	---	2.7	U	NP
Arsenic	0.05	U	---	0.42	UW	---	0.42	U	NP	0.20	U	---	0.20	U	NP	0.48	U	---	0.46	U	NP
Barium	0.40	B	---	515.00		---	519.00		NP	0.55	B	J	0.47	B	NP	0.12	B	J	0.17	B	NP
Beryllium	0.04	U	---	0.06	B	J	0.05	B	NP	0.16	U	---	0.21	B	NP	0.05	B	U	0.05	B	NP
Cadmium	0.12	U	---	0.11	U	---	0.12	U	NP	0.47	U	---	0.51	U	NP	0.12	U	---	0.13	U	NP
Calcium	250		---	218000		---	229000		NP	299	B	J	271	B	NP	96.5	B	J	117	B	NP
Chromium	0.22	B	---	3.40		---	3.30		NP	0.71	B	J	0.66	U	NP	0.16	B	J	0.24	B	NP
Cobalt	0.27	U	---	0.24	U	---	0.26	U	NP	1.00	U	---	1.10	U	NP	0.28	U	---	0.29	U	NP
Copper	1.40		---	0.15	U	---	0.16	U	NP	5.10		---	2.20	B	NP	49.50		---	45.30		NP
Iron	29.0		---	2.8	B	U	3.8	B	NP	15.5	B	U	33.9		NP	217.0		---	170.0		NP
Lead	1.28		---	0.52		---	0.45		NP	0.40	B	J	0.17	U	NP	0.05	UW	UJ	0.05	U	NP
Magnesium	268.0		---	3230.0		---	3270.0		NP	77.5	B	J	86.2	B	NP	162.0	B	J	178.0	B	NP
Manganese	0.76		---	1.10		---	1.40		NP	0.21	U	---	0.23	U	NP	4.00		---	3.90		NP
Mercury	0.04	B	---	0.03	U	---	0.04	B	NP	0.07	B	U	0.06	B	NP	0.08	B	U	0.05	B	NP
Nickel	0.90	U	---	0.81	U	---	0.86	U	NP	3.50	U	---	3.80	U	NP	0.92	U	---	0.95	U	NP
Potassium	3120		---	316		---	387		NP	581	B	J	537	B	NP	2660		---	2660		NP
Selenium	0.49	U	---	0.42	U	UJ	0.42	U	NP	0.20	U	---	0.20	U	NP	0.49	BW	J	0.36	B	NP
Silver	0.16	U	---	0.14	U	---	0.15	U	NP	0.61	U	---	0.66	U	NP	0.16	U	---	0.17	U	NP
Sodium	715.0		---	5210.0		---	5380.0		NP	394.0	B	J	508.0	B	NP	834.0		---	1034.0		NP
Thallium	0.05	U	---	0.04	UW	UJ	0.04	U	NP	0.20	U	---	0.20	U	NP	0.05	U	---	0.05	U	NP
Vanadium	0.16	U	---	0.15	U	---	0.16	U	NP	0.63	U	---	0.68	U	NP	0.17	U	---	0.17	U	NP
Zinc	42.6		---	50.7		---	56.0		NP	4.1		---	6.0		NP	31.9		---	33.5		NP

* Data Qualifier. ** Data Validation.

Table A-8.1 Metals

Compound (mg/Kg)	1D-LM-M-I	Q*	DV**	1D-LM-M-I-D	Q*	DV**	1E-LM-B-I	Q*	DV**	1E-LM-F-I	Q*	DV**	1E-LM-L-I	Q*	DV**	1E-LM-M-I	Q*	DV**	1F-LM-B-I	Q*	DV**
Aluminum	1.4	B	J	1.8	B	NP	5.5	B	J	6.8	B	J	1.5	B	J	1.1	B	J	6.6	B	
Antimony	2.6	U	---	2.7	U	NP	3.3		---	11.1	U	---	2.4	U	---	2.4	U	---	2.7	U	---
Arsenic	0.47	U	---	0.46	U	NP	0.50	UW	---	0.19	U	J	0.49	U	---	0.40	U	---	1.40	BW	J
Barium	0.21	B	J	0.29	B	NP	540.00		---	0.75	B	U	0.21	B	J	0.72	B	J	565.00		---
Beryllium	0.04	U	---	0.05	B	NP	0.05	B	J	0.22	B	---	0.04	U	---	0.05	B	U	0.04	U	---
Cadmium	0.13	U	---	0.13	U	NP	0.13	U	---	0.54	U	---	0.39		---	0.12	U	---	0.13	U	---
Calcium	133	B	J	157	B	NP	227000		---	338	B	J	132	B	J	336		---	207000		---
Chromium	0.16	U	---	0.24	B	NP	3.30		---	0.70	U	---	0.15	U	---	0.15	U	---	3.40		---
Cobalt	0.28	U	---	0.29	U	NP	0.30	U	---	1.20	U	---	0.25	U	---	0.26	U	---	0.29	B	J
Copper	2.10		---	1.50		NP	0.18	U	---	3.40	B	J	40.30		---	1.50		---	0.18	U	---
Iron	27.9		U	22.5		NP	4.5	B	J	21.7		U	204.0		---	25.2		U	4.3	B	J
Lead	0.10	BW	J	0.05	U	NP	0.32		---	0.19	U	---	0.05	U	---	0.04	U	---	0.95		---
Magnesium	277.0		---	199.0	B	NP	3230.0		---	91.9	B	J	168.0	B	J	283.0		---	3050.0		---
Manganese	0.17	B	J	0.21	B	NP	1.40		---	0.24	U	---	4.40		---	0.18	B	J	1.60		---
Mercury	0.06	B	U	0.04	B	NP	0.04	B	U	0.05	U	---	0.20		U	0.05	B	U	0.04	U	---
Nickel	0.93	U	---	0.96	U	NP	1.00	U	---	4.00	U	---	0.85	U	---	0.86	U	---	0.98	U	---
Potassium	3560		---	2600		NP	385		---	870	B	J	2380		---	3710		---	274		U
Selenium	0.24	UW	---	0.23	U	NP	0.66	BE	J	0.19	UW	UJ	0.42	BW	J	0.20	UW	UJ	0.57	BW	J
Silver	0.16	U	UJ	0.17	U	NP	0.17	U	---	0.70	U	---	0.15	U	---	0.15	U	---	0.17	U	---
Sodium	573.0		---	505.0		NP	5250.0		---	711.0	B	J	1050.0		---	507.0		---	5590.0		---
Thallium	0.05	U	---	0.05	U	NP	0.05	UW	UJ	0.19	U	---	0.05	U	---	0.04	U	---	0.04	UW	UJ
Vanadium	0.17	U	---	0.17	U	NP	0.18	U	---	0.72	U	---	0.15	U	---	0.16	U	---	0.18	U	---
Zinc	33.6		---	26.6		NP	70.4		---	5.4		---	36.6		---	30.6		---	70.1		---

* Data Qualifier. ** Data Validation.

Compound (mg/Kg)	Table A-8.1 Metals																							
	1F-LM-F-I	Q*	DV**	1F-LM-L-I	Q*	DV**	1F-LM-M-I	Q*	DV**	1G-LM-B-I	Q*	DV**	1G-LM-F-I	Q*	DV**	1G-LM-L-I	Q*	DV**	1G-LM-M-I	Q*	DV**
Aluminum	2.1	B	U	3.7	B	U	1.8	B	U	5.1	B	U	4.9	B	U	3.6	B	U	1.4	B	J
Antimony	2.7	U	---	4.2	U	---	2.6	U	---	2.7	U	---	2.6	U	---	5.5	U	---	2.1	U	---
Arsenic	0.05	U	---	0.07	U	---	0.05	U	---	0.62	BW	J	0.05	UW	UJ	0.10	U	---	0.05	UW	---
Barium	0.63	B	J	0.44	B	J	0.32	B	J	560.00		---	1.80	B	J	0.41	B	J	0.37	B	J
Beryllium	0.04	U	---	0.07	U	---	0.04	U	---	0.04	U	---	0.04	U	---	0.09	U	---	0.03	U	---
Cadmium	0.13	U	---	0.20	U	---	0.13	U	---	0.13	U	---	0.12	U	---	0.27	U	---	0.10	U	---
Calcium	326		---	271	B	J	169	B	J	204000		---	665		---	192	B	J	221		---
Chromium	0.28	B	J	0.32	B	J	0.55		---	2.90		---	0.26	B	J	0.77	B	J	0.39		--
Cobalt	0.29	U	---	0.45	U	---	0.29	U	---	0.29	U	---	0.28	U	---	0.60	U	---	0.22	U	---
Copper	0.28	B	J	64.80		---	2.20		---	0.18	U	---	0.85	B	U	55.30		---	2.00		---
Iron	4.0	B	J	137.0		---	34.5		---	3.7	B	J	22.3		---	258.0		---	35.9		---
Lead	0.05	U	---	0.06	UW	UJ	0.05	B	UJ	1.80		---	0.12	B	U	0.12	BW	UJ	0.08	B	UJ
Magnesium	25.4	B	J	204.0	B	J	244.0		---	3190.0		---	73.8	B	J	182.0	B	J	275.0		---
Manganese	0.09	B	J	5.90		---	0.29	B	J	1.60		---	0.73		---	5.40		---	0.59		---
Mercury	0.05	U	---	0.04	B	J	0.03	U	---	0.03	U	---	0.05	U	---	0.19		---	0.03	U	---
Nickel	0.97	U	---	1.50	U	---	0.95	U	---	0.98	U	---	0.92	U	---	2.00	U	---	0.75	U	---
Potassium	182	B	U	2580		---	2880		---	180	B	U	718		U	2800		---	3570		---
Selenium	0.48	U	---	0.77	B	J	0.46	UW	UJ	0.43	U	---	0.46	U	---	0.99	U	---	0.48	UW	UJ
Silver	0.17	U	---	0.26	U	---	0.17	U	---	0.17	U	---	0.16	U	---	0.35	U	---	0.13	U	---
Sodium	161.0	B	J	807.0		---	619.0		---	5230.0		---	1050.0		---	1010.0		---	548.0		---
Thallium	0.05	U	---	0.07	U	---	0.05	U	---	0.04	U	---	0.05	U	---	0.10	U	---	0.05	U	---
Vanadium	0.17	U	---	0.27	U	---	0.17	U	---	0.18	U	---	0.17	U	---	0.36	U	---	0.13	U	---
Zinc	2.4		---	47.7		---	39.8		---	67.4		---	6.7		---	43.0		---	37.5		---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1A-LM-B-I	Q*	DV**	1A-LM-L-I	Q*	DV**	1A-LM-M-I	Q*	DV**	1B-LM-B-I	Q*	DV**	1B-LM-F-I	Q*	DV**	1B-LM-L-I	Q*	DV**	1B-LM-M-I	Q*	DV**
Alpha-BHC	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Beta-BHC	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Delta-BHC	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Gamma-BHC	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Heptachlor	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Aldrin	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Heptachlor epoxide	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Endosulfan I	46	U	---	48	U	---	46	U	---	46	U	---	48	U	---	48	U	---	48	U	---
Dieldrin	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
4,4'-DDE	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
Endrin	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
Endosulfan II	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
4,4'-DDD	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
Endosulfan sulfate	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
4,4'-DDT	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
Methoxychlor	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---
Endrin ketone	93	U	---	96	U	---	92	U	---	92	U	---	95	U	---	95	U	---	95	U	---
Alpha-chlordane	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---
Gamma-chlordane	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---
Toxaphene	930	U	---	960	U	---	920	U	---	920	U	---	950	U	---	950	U	---	950	U	---
Aroclor-1016	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---
Aroclor-1221	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---

* Data Qualifier. ** Data Validation.

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1A-LM-B-I	Q*	DV**	1A-LM-L-I	Q*	DV**	1A-LM-M-I	Q*	DV**	1B-LM-B-I	Q*	DV**	1B-LM-F-I	Q*	DV**	1B-LM-L-I	Q*	DV**	1B-LM-M-I	Q*	DV**
Aroclor-1232	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---
Aroclor-1242	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---
Aroclor-1248	460	U	---	480	U	---	460	U	---	460	U	---	480	U	---	480	U	---	480	U	---
Aroclor-1254	930	U	---	960	U	---	920	U	---	920	U	---	950	U	---	950	U	---	950	U	---
Aroclor-1260	930	U	---	960	U	---	920	U	---	920	U	---	950	U	---	950	U	---	950	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1C-LM-B-I	Q*	DV**	1C-LM-B-I-D	Q*	DV**	1C-LM-F-I	Q*	DV**	1C-LM-F-I-D	Q*	DV**	1C-LM-L-I	Q*	DV**	1C-LM-L-I-D	Q*	DV**	1C-LM-M-I	Q*	DV**
Alpha-BHC	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Beta-BHC	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Delta-BHC	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Gamma-BHC	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Heptachlor	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Aldrin	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Heptachlor epoxide	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Endosulfan I	47	U	---	47	U	---	48	U	---	47	U	---	47	U	---	47	U	---	48	U	---
Dieldrin	94	U	---	95	U	---	95	U	---	95	U	.	93	U	---	95	U	---	95	U	---
4,4'-DDE	94	U	---	95	U	---	95	U	---	95	U	---	93	U	---	95	U	---	95	U	---
Endrin	94	U	---	95	U	---	95	U	---	95	U	---	93	U	---	95	U	---	95	U	---
Endosulfan II	94	U	---	95	U	---	95	U	---	95	U	---	93	U	---	95	U	---	95	U	---
4,4'-DDD	94	U	---	95	U	---	95	U	---	95	U	---	93	U	---	95	U	---	95	U	---
Endosulfan sulfate	94	U	---	95	U	---	95	U	---	95	U	---	93	U	---	95	U	---	95	U	---
4,4'-DDT	94	U	---	95	U	---	95	U	---	95	U	---	93	U	---	95	U	---	95	U	---
Methoxychlor	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---
Endrin ketone	94	U	---	95	U	---	95	U	---	95	U	---	93	U	---	95	U	---	95	U	---
Alpha-chlordane	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---
Gamma-chlordane	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---
Toxaphene	940	U	---	950	U	---	950	U	---	950	U	---	930	U	---	950	U	---	950	U	---
Aroclor-1016	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---
Aroclor-1221	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1C-LM-B-I	Q*	DV**	1C-LM-B-I-D	Q*	DV**	1C-LM-F-I	Q*	DV**	1C-LM-F-I-D	Q*	DV**	1C-LM-L-I	Q*	DV**	1C-LM-L-I-D	Q*	DV**	1C-LM-M-I	Q*	DV**
Aroclor-1232	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---
Aroclor-1242	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---
Aroclor-1248	470	U	---	470	U	---	480	U	---	470	U	---	470	U	---	470	U	---	480	U	---
Aroclor-1254	940	U	---	950	U	---	950	U	---	950	U	---	930	U	---	950	U	---	950	U	---
Aroclor-1260	940	U	---	950	U	---	950	U	---	950	U	---	930	U	---	950	U	---	950	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1C-LM-M-I-D	Q*	DV**	1D-LM-B-I	Q*	DV**	1D-LM-B-I-D	Q*	DV**	1D-LM-F-I	Q*	DV**	1D-LM-F-I-D	Q*	DV**	1D-LM-L-I	Q*	DV**	1D-LM-L-I-D	Q*	DV**
Alpha-BHC	48	U	---	47	U	---	48	U	---	48	U	---	47	U	---	46	U	---	48	U	---
Beta-BHC	48	U	---	47	U	---	48	U	---	48	U	---	47	U	---	46	U	---	48	U	---
Delta-BHC	48	U	---	47	U	---	48	U	---	48	U	---	47	U	---	46	U	---	48	U	---
Gamma-BHC	48	U	---	47	U	---	48	U	---	48	U	---	47	U	---	46	U	---	48	U	---
Heptachlor	48	U	---	47	U	---	48	U	---	48	U	---	47	U	---	46	U	---	48	U	---
Aldrin	48	U	---	47	U	---	48	U	---	48	U	---	47	U	---	46	U	---	48	U	---
Heptachlor epoxide	48	U	---	47	U	---	48	U	---	48	U	---	47	U	---	46	U	---	48	U	---
Endosulfan I	48	U	---	47	U	---	48	U	---	180	UX	---	47	U	---	46	U	---	48	U	---
Dieldrin	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
4,4'-DDE	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
Endrin	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
Endosulfan II	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
4,4'-DDD	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
Endosulfan sulfate	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
4,4'-DDT	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
Methoxychlor	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---
Endrin ketone	95	U	---	93	U	---	95	U	---	95	U	---	93	U	---	91	U	---	95	U	---
Alpha-chlordane	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---
Gamma-chlordane	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---
Toxaphene	950	U	---	930	U	---	950	U	---	950	U	---	930	U	---	910	U	---	950	U	---
Aroclor-1016	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---
Aroclor-1221	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1C-LM-M-I-D	Q*	DV**	1D-LM-B-I	Q*	DV**	1D-LM-B-I-D	Q*	DV**	1D-LM-F-I	Q*	DV**	1D-LM-F-I-D	Q*	DV**	1D-LM-L-I	Q*	DV**	1D-LM-L-I-D	Q*	DV**
Aroclor-1232	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---
Aroclor-1242	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---
Aroclor-1248	480	U	---	470	U	---	480	U	---	480	U	---	470	U	---	460	U	---	480	U	---
Aroclor-1254	950	U	---	930	U	---	950	U	---	950	U	---	930	U	---	910	U	---	950	U	---
Aroclor-1260	950	U	---	930	U	---	950	U	---	950	U	---	930	U	---	910	U	---	950	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1D-LM-M-I	Q*	DV**	1D-LM-M-I-D	Q*	DV**	1E-LM-B-I	Q*	DV**	1E-LM-F-I	Q*	DV**	1E-LM-L-I	Q*	DV**	1E-LM-M-I	Q*	DV**	1F-LM-B-I	Q*	DV**
Alpha-BHC	48	U	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	--	46	U	---
Beta-BHC	48	U	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	---	46	U	---
Delta-BHC	90	UX	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	---	46	U	---
Gamma-BHC	48	U	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	---	46	U	---
Heptachlor	48	U	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	---	46	U	---
Aldrin	48	U	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	---	46	U	---
Heptachlor epoxide	60	UX	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	---	46	U	---
Endosulfan I	48	U	---	47	U	---	47	U	J	47	U	---	46	U	---	47	U	---	46	U	---
Dieldrin	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
4,4'-DDE	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
Endrin	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
Endosulfan II	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
4,4'-DDD	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
Endosulfan sulfate	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
4,4'-DDT	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
Methoxychlor	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U		460	U	---
Endrin ketone	95	U	---	93	U	---	95	U	J	95	U	---	92	U	---	93	U	---	92	U	---
Alpha-chlordane	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U	---	460	U	---
Gamma-chlordane	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U	---	460	U	---
Toxaphene	950	U	---	930	U	---	950	U	J	950	U	---	920	U	---	930	U	---	920	U	---
Aroclor-1016	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U	---	460	U	---
Aroclor-1221	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U	---	460	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1D-LM-M-I	Q*	DV**	1D-LM-M-I-D	Q*	DV**	1E-LM-B-I	Q*	DV**	1E-LM-F-I	Q*	DV**	1E-LM-L-I	Q*	DV**	1E-LM-M-I	Q*	DV**	1F-LM-B-I	Q*	DV**
Aroclor-1232	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U	---	460	U	---
Aroclor-1242	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U	---	460	U	---
Aroclor-1248	480	U	---	470	U	---	470	U	J	470	U	---	460	U	---	470	U	---	460	U	---
Aroclor-1254	950	U	---	930	U	---	950	U	J	950	U	---	920	U	---	930	U	---	920	U	---
Aroclor-1260	950	U	---	930	U	---	950	U	J	950	U	---	920	U	---	930	U	---	920	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1F-LM-F-I	Q*	DV**	1F-LM-L-I	Q*	DV**	1F-LM-M-I	Q*	DV**	1G-LM-B-I	Q*	DV**	1G-LM-F-I	Q*	DV**	1G-LM-L-I	Q*	DV**	1G-LM-M-I	Q*	DV**
Alpha-BHC	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Beta-BHC	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Delta-BHC	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Gamma-BHC	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Heptachlor	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Aldrin	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Heptachlor epoxide	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Endosulfan I	48	U	---	47	U	---	46	U	---	47	U	---	48	U	---	47	U	---	47	U	---
Dieldrin	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
4,4'-DDE	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
Endrin	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
Endosulfan II	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
4,4'-DDD	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
Endosulfan sulfate	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
4,4'-DDT	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
Methoxychlor	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Endrin ketone	95	U	---	94	U	---	91	U	---	93	U	---	95	U	---	95	U	---	94	U	---
Alpha-chlordane	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Gamma-chlordane	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Toxaphene	950	U	---	940	U	---	910	U	---	930	U	---	950	U	---	950	U	---	940	U	---
Aroclor-1016	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1221	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---

Compound (ug/Kg)	Table A-8.2 Pesticides/PCB's																							
	1F-LM-F-I	Q*	DV**	1F-LM-L-I	Q*	DV**	1F-LM-M-I	Q*	DV**	1G-LM-B-I	Q*	DV**	1G-LM-F-I	Q*	DV**	1G-LM-L-I	Q*	DV**	1G-LM-M-I	Q*	DV**
Aroclor-1232	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1242	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1248	480	U	---	470	U	---	460	U	---	470	U	---	480	U	---	470	U	---	470	U	---
Aroclor-1254	950	U	---	940	U	---	910	U	---	930	U	---	950	U	---	950	U	---	940	U	---
Aroclor-1260	950	U	---	940	U	---	910	U	---	930	U	---	950	U	---	950	U	---	940	U	---

Parameter (pCi/g)	Table A-8.3 Gross Alpha/Beta																			
	1A-LM-B-I	DV**	1A-LM-L-I	DV**	1A-LM-M-I	DV**	1B-LM-B-I	DV**	1B-LM-F-I	DV**	1B-LM-L-I	DV**	1B-LM-M-I	DV**	1C-LM-B-I	DV**	1C-LM-B-I-D	DV**
Alpha	7.80	J	ND	---	ND	---	ND	---	0.58	J	0.22	J	0.24	J	ND	---	ND	---
Sigma Error +/-	5.73	---	-	---	-	---	-	---	0.13	---	0.11	---	0.13	---	-	---	-	---
MDA*	7.64	---	0.34	UJ	0.24	UJ	10.48	UJ	0.10	---	0.10	---	0.14	---	9.75	UJ	12.72	UJ
Beta	30.60	---	2.29	---	2.13	---	21.05	---	0.41	---	2.68	---	2.34	---	28.20	---	37.64	---
Sigma Error +/-	4.92	---	0.20	---	0.15	---	4.86	---	0.08	---	0.16	---	0.16	---	4.52	---	5.12	---
MDA*	5.76	---	0.17	---	0.11	---	6.68	---	0.10	---	0.11	---	0.12	---	5.11	---	5.46	---

Parameter (pCi/g)	Table A-8.3 Gross Alpha/Beta																	
	1C-LM-F-I	DV**	1C-LM-F-I-D	DV**	1C-LM-L-I	DV**	1C-LM-L-I-D	DV**	1C-LM-M-I	DV**	1C-LM-M-I-D	DV**	1D-LM-B-I	DV**	1D-LM-B-I-D	DV**	1D-LM-F-I	DV**
Alpha	0.02	---	1.42	J	0.34	J	ND	---	0.10	---	0.33	J	7.29	---	6.02	---	0.03	---
Sigma Error +/-	0.03	---	0.16	---	0.16	---	-	---	0.11	---	0.18	---	5.35	---	4.73	---	0.03	---
MDA*	0.05	UJ	0.06	---	0.17	---	0.39	UJ	0.17	UJ	0.24	---	6.78	---	6.22	U	0.05	U
Beta	0.11	---	0.81	---	3.06	---	2.57	---	2.30	---	2.23	---	32.82	---	32.71	---	0.13	---
Sigma Error +/-	0.06	---	0.10	---	0.20	---	0.18	---	0.16	---	0.17	---	5.44	---	5.61	---	0.06	---
MDA*	0.10	---	0.11	---	0.13	---	0.13	---	0.12	---	0.14	---	6.57	---	5.89	---	0.10	---

* Minimum Detectable Activity. ** Data Validation.

Parameter (pCi/g)	1D-LM-F-I-D	DV**	1D-LM-L-I	DV**	1D-LM-L-I-D	DV**	1D-LM-M-I	DV**	1D-LM-M-I-D	DV**	1E-LM-B-I	DV**	1E-LM-F-I	DV**	1E-LM-L-I	DV**	1E-LM-M-I	DV**
Table A-8.3 Gross Alpha/Beta																		
Alpha	0.04	---	0.36	---	0.56	---	0.09	---	0.15	---	ND	---	0.11	---	0.23	---	0.16	---
Sigma Error +/-	0.06	---	0.16	---	0.23	---	0.11	---	0.13	---	-	---	0.08	---	0.15	---	0.16	---
MDA*	0.10	U	0.19	---	0.27	---	0.18	U	0.19	U	12.21	U	0.11	U	0.21	---	0.26	U
Beta	0.24	---	2.27	---	2.37	---	2.02	---	1.95	---	37.43	---	0.16	---	2.75	---	2.59	---
Sigma Error +/-	0.07	---	0.15	---	0.17	---	0.14	---	0.14	---	6.36	---	0.08	---	0.17	---	0.19	---
MDA*	0.09	---	0.10	---	0.11	---	0.09	---	0.12	---	7.59	---	0.13	---	0.12	---	0.15	---

Table A-8.3 Gross Alpha/Beta																				
Parameter (pCi/g)	1F-LM-B-I	DV**	1F-LM-F-I	DV**	1F-LM-L-I	DV**	1F-LM-M-I	DV**	1G-LM-B-I	DV**	1G-LM-F-I	DV**	1G-LM-L-I	DV**	1G-LM-M-I	DV**
Alpha	ND	---	ND	---	0.40	J	0.26	J	ND	---	0.09	---	0.34	J	0.13	---
Sigma Error +/-	-	---	-	---	0.17	---	0.14	---	-	---	0.14	---	0.20	---	0.11	---
MDA*	12.12	UJ	0.05	UJ	0.19	---	0.17	---	12.20	UJ	0.23	UJ	0.29	---	0.16	UJ
Beta	36.09	---	ND	---	2.67	---	2.46	---	29.72	---	0.32	---	2.58	---	3.07	---
Sigma Error +/-	6.01	---	-	---	0.18	---	0.17	---	4.78	---	0.08	---	0.18	---	0.19	---
MDA*	7.11	---	0.10	U	0.13	---	0.11	---	5.48	---	0.11	---	0.12	---	0.12	---

Table A-8.4 Gamma Spectroscopy

Nuclide (pCi/g)	1A-LM-B-I	DV**	1A-LM-F-I	DV**	1A-LM-L-I	DV**	1A-LM-M-I	DV**	1B-LM-B-I	DV**	1B-LM-F-I	DV**	1B-LM-L-I	DV**	1B-LM-M-I	DV**	1C-LM-B-I	DV**
Cesium 137	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	0.32	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.04	---	-	---
MDA*	0.20	U	0.40	U	0.20	U	0.20	U	0.10	U	0.20	U	0.20	U	0.10	---	0.10	U
Potassium 40	ND	---	ND	---	ND	---	2.70	---	ND	---	ND	---	10.20	---	4.10	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	0.40	---	-	---	-	---	0.70	---	0.50	---	-	---
MDA*	0.50	6.0U	11.10	U	3.10	U	2.50	---	2.60	U	5.20	U	0.80	---	1.90	---	1.70	U
Cadmium 109	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Lead 212	ND	---	ND	---	ND	---	ND	---	0.14	---	ND	---	ND	---	ND	---	0.14	---
Sigma Error +/-	-	---	-	---	-	---	-	---	0.03	---	-	---	-	---	-	---	0.04	---
MDA*	0.30	U	0.60	U	0.20	U	0.20	U	0.10	---	0.30	U	0.30	U	0.10	U	0.10	---
Radium 226	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	0.70	U	0.30	U	0.20	U	0.20	U	0.30	U	0.30	U	0.20	U	0.20	U
Bismuth 214	0.40	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	0.10	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	---	0.70	U	0.40	U	0.20	U	0.30	U	0.40	U	0.40	U	0.30	U	0.20	U
Lead 210	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Uranium 238	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-8.4 Gamma Spectroscopy																			
	1A-LM-B-I	DV**	1A-LM-F-I	DV**	1A-LM-L-I	DV**	1A-LM-M-I	DV**	1B-LM-B-I	DV**	1B-LM-F-I	DV**	1B-LM-L-I	DV**	1B-LM-M-I	DV**	1C-LM-B-I	DV**
Radium 228	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Thallium 208	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

Nuclide (pCi/g)	Table A-8.4 Gamma Spectroscopy																			
	1C-LM-B-I-D	DV**	1C-LM-F-I	DV**	1C-LM-F-I-D	DV**	1C-LM-L-I	DV**	1C-LM-L-I-D	DV**	1C-LM-M-I	DV**	1C-LM-M-I-D	DV**	1D-LM-B-I	DV**	1D-LM-B-I-D	DV**
Cesium 137	ND	---	ND	---	ND	---	0.14	---	ND	---	0.32	---	0.75	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	0.03	---	-	---	0.04	---	0.06	---	-	---	-	---
MDA*	0.10	U	0.10	U	0.20	U	0.10	---	0.20	U	0.10	---	0.10	---	0.20	U	0.10	U
Potassium 40	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	2.60	U	1.70	U	5.10	U	1.80	U	4.50	U	2.40	U	4.40	U	3.60	U	2.30	U
Cadmium 109	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	2.00	U	0.90	U
Lead 212	ND	---	ND	---	ND	---	ND	---	0.38	---	ND	---	ND	---	ND	---	0.13	---
Sigma Error +/-	-	---	-	---	-	---	-	---	0.04	---	-	---	-	---	-	---	0.01	---
MDA*	0.20	U	0.20	U	0.30	U	0.10	U	0.20	---	0.20	U	0.30	U	0.30	U	0.10	---
Radium 226	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.20	U	0.30	U	0.10	U	0.20	U	0.20	U	0.40	U	-	---	-	---
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	0.20	U	0.30	U	0.20	U	0.30	U	0.20	U	0.50	U	0.40	U	0.30	U
Lead 210	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	1.80	U	0.05	U
Uranium 238	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	1.70	U	0.90	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-8.4 Gamma Spectroscopy																	
	1C-LM-B-I-D	DV**	1C-LM-F-I	DV**	1C-LM-F-I-D	DV**	1C-LM-L-I	DV**	1C-LM-L-I-D	DV**	1C-LM-M-I	DV**	1C-LM-M-I-D	DV**	1D-LM-B-I	DV**	1D-LM-B-I-D	DV**
Radium 228	-	---	-	---	-	---	-	---	-	---	-	---	-	---	1.40	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.20	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.80	---	0.30	U
Thallium 208	-	---	-	---	-	---	-	---	-	---	-	---	-	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.20	U	0.10	U

* Minimum Detectable Activity. ** Data Validation.

Table A-8.4 Gamma Spectroscopy

Nuclide (pCi/g)	1D-LM-F-I	DV**	1D-LM-F-I-D	DV**	1D-LM-L-I	DV**	1D-LM-L-I-D	DV**	1D-LM-M-I	DV**	1D-LM-M-I-D	DV**	1E-LM-B-I	DV**	1E-LM-F-I	DV**	1E-LM-L-I	DV**
Cesium 137	ND	---	ND	---	0.22	---	0.24	---	0.43	---	0.32	---	ND	---	ND	---	0.24	---
Sigma Error +/-	-	---	-	---	0.03	---	0.03	---	0.04	---	0.03	---	-	---	-	---	0.04	---
MDA*	0.10	U	0.10	U	0.10	---	0.10	---	0.10	---	0.10	---	0.10	U	0.10	U	0.10	---
Potassium 40	ND	---	ND	---	ND	---	4.70	---	3.90	---	ND	---	ND	---	ND	---	3.20	---
Sigma Error +/-	-	---	-	---	-	---	0.40	---	0.30	---	-	---	-	---	-	---	0.40	---
MDA*	1.70	U	2.40	U	4.60	U	0.80	---	2.50	---	4.70	U	2.40	U	2.90	U	1.90	---
Cadmium 109	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.80	U	1.00	U	1.50	U	1.00	U	1.30	U	1.40	U	1.30	U	1.10	U	1.00	U
Lead 212	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	0.50	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	0.10	---	-	---	-	---
MDA*	0.20	U	0.20	U	0.20	U	0.20	U	0.20	U	0.20	U	0.20	---	0.20	U	0.20	U
Radium 226	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.20	U	0.20	U	0.30	U	0.30	U	0.30	U	0.20	U	0.30	U	0.30	U
Lead 210	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.10	U	1.20	U	1.40	U	0.10	U	1.00	U	1.30	U	1.50	U	0.10	U	0.07	U
Uranium 238	ND	---	ND	---	ND	---	3.80	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	0.20	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.30	U	1.00	U	1.20	U	0.70	---	1.00	U	1.10	U	1.20	U	1.10	U	1.10	U

* Minimum Detectable Activity. ** Data Validation.

Nuclide (pCi/g)	Table A-8.4 Gamma Spectroscopy																	
	1D-LM-F-I	DV**	1D-LM-F-I-D	DV**	1D-LM-L-I	DV**	1D-LM-L-I-D	DV**	1D-LM-M-I	DV**	1D-LM-M-I-D	DV**	1E-LM-B-I	DV**	1E-LM-F-I	DV**	1E-LM-L-I	DV**
Radium 228	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.40	U	0.50	U	0.50	U	0.50	U	0.50	U	0.50	U	0.70	U	0.60	U	0.50	U
Thallium 208	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	0.30	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.10	---	-	---
MDA*	0.10	U	0.10	U	0.10	U	0.10	U	0.20	U	0.10	U	0.10	U	0.10	---	0.10	U

Table A-8.4 Gamma Spectroscopy

Nuclide (pCi/g)	1E-LM-M-I	DV**	1F-LM-B-I	DV**	1F-LM-F-I	DV**	1F-LM-L-I	DV**	1F-LM-M-I	DV**	1G-LM-B-I	DV**	1G-LM-F-I	DV**	1G-LM-L-I	DV**	1G-LM-M-I	DV**
Cesium 137	ND	---	ND	---	ND	---	0.21	---	0.77	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	0.02	---	0.07	---	-	---	-	---	-	---	-	---
MDA*	0.10	U	0.10	U	0.10	U	0.10	---	0.10	---	0.10	U	0.20	U	0.20	U	0.10	U
Potassium 40	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	3.70	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.50	---
MDA*	5.20	U	2.20	U	1.30	U	3.70	U	3.70	U	1.80	U	5.90	U	6.50	U	2.20	---
Cadmium 109	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.50	U	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Lead 212	ND	---	0.20	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	0.03	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.20	U	0.10	---	0.10	U	0.20	U	0.30	U	0.10	U	0.40	U	0.30	U	0.10	U
Radium 226	-	---	ND	---	ND	---	ND	---	ND	---	0.26	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	0.03	---	-	---	-	---	-	---
MDA*	-	---	0.20	U	0.10	U	0.20	U	0.30	U	0.20	---	0.40	U	0.40	U	0.20	U
Bismuth 214	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.30	U	0.30	U	0.20	U	0.20	U	0.40	U	0.20	U	0.50	U	0.50	U	0.20	U
Lead 210	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.40	U	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Uranium 238	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	1.30	U	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

Nuclide (pCi/g)	Table A-8.4 Gamma Spectroscopy																			
	1E-LM-M-I	DV**	1F-LM-B-I	DV**	1F-LM-F-I	DV**	1F-LM-L-I	DV**	1F-LM-M-I	DV**	1G-LM-B-I	DV**	1G-LM-F-I	DV**	1G-LM-L-I	DV**	1G-LM-M-I	DV**
Radium 228	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.60	u	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Thallium 208	ND	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---
MDA*	0.10	u	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---

Table A-8.5 Tritium

Nuclide (pCi/g)	1A-LM-B-I	DV**	1A-LM-F-I	DV**	1A-LM-L-I	DV**	1A-LM-M-I	DV**	1B-LM-B-I	DV**	1B-LM-F-I	DV**	1B-LM-L-I	DV**	1B-LM-M-I	DV**	1C-LM-B-I	DV**
Tritium	ND	---	12.88	---	0.10	---	ND	---	ND	---	0.14	---	ND	---	ND	---	ND	---
Sigma Error +/-	-	---	1.29	---	0.05	---	-	---	-	---	0.03	---	-	---	-	---	-	---
MDA*	0.05	U	0.13	---	0.11	U	0.13	U	0.05	U	0.05	---	0.08	U	0.12	U	0.05	U

Nuclide (pCi/g)	1C-LM-B-I-D	DV**	1C-LM-F-I	DV**	1C-LM-F-I-D	DV**	1C-LM-L-I	DV**	1C-LM-L-I-D	DV**	1C-LM-M-I	DV**	1C-LM-M-I-D	DV**	1D-LM-B-I	DV**	1D-LM-B-I-D	DV**
Table A-8.5 Tritium																		
Tritium	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	ND	---	0.54	---
Sigma Error +/-	-	---	-	---	-	---	-	---	-	---	-	---	-	---	-	---	0.07	---
MDA*	0.10	U	0.05	U	0.05	U	0.06	U	0.09	U	0.13	U	0.12	U	0.09	U	0.09	---

Table A-8.5 Tritium

Nuclide (pCi/g)	1D-LM-F-I	DV**	1D-LM-F-I-D	DV**	1D-LM-L-I	DV**	1D-LM-L-I-D	DV**	1D-LM-M-I	DV**	1D-LM-M-I-D	DV**	1E-LM-B-I	DV**	1E-LM-F-I	DV**	1E-LM-L-I	DV**
Tritium	ND	---	ND	---	ND	---	25.55	---	ND	---	ND	---	ND	---	84.59	---	6.60	---
Sigma Error +/-	-	---	-	---	-	---	2.52	---	-	---	-	---	-	---	8.30	---	0.67	---
MDA*	0.10	U	0.09	U	0.21	U	0.16	---	0.20	U	0.19	U	0.09	U	0.10	---	0.17	---

Nuclide (pCi/g)	Table A-8.5 Tritium																			
	1E-LM-M-I	DV**	1F-LM-B-I	DV**	1F-LM-F-I	DV**	1F-LM-L-I	DV**	1F-LM-M-I	DV**	1G-LM-B-I	DV**	1G-LM-F-I	DV**	1G-LM-L-I	DV**	1G-LM-M-I	DV**
Tritium	0.36	---	ND	---	ND	---	0.11	---	ND	---	ND	---	28.41	---	ND	---	ND	---
Sigma Error +/-	0.11	---	-	---	-	---	0.05	---	-	---	-	---	2.80	---	-	---	-	---
MDA*	0.23	---	0.05	U	0.05	U	0.11	U	0.13	U	0.05	U	0.08	---	0.12	U	0.11	U

Table A-9.1
RBP-I Habitat Assessment Data

Habitat Parameter (RANGE OF VALUES)	TDTS STATION IDENTIFICATIONS								
	HMC-1[1]	HMC-2	HMC-3	HMC-4	HMC-5	GRC-1[1]	GRC-2	HHC-1	HHC-2
Bottom Substrate (0-20)	11	9	9	11	8	10	9	11	6
Stream Flow and/or Stream Velocity (0-20)	9	11	11	18	13	16	11	9	2
Channel Aeration (0-15)	11	11	11	7	11	9	11	11	15
Bottom Scouring and Deposition (0-15)	12	12	12	9	10	9	12	12	7
Pool/Riffle or Run/Bend Ration (0-15)	8	10	11	13	12	12	11	8	2
Bank Stability (0-10)	10	10	10	9	8	9	10	10	10
Bank Vegetative Survey (0-10)	10	9	10	8	10	9	10	10	10
Streamside Cover (0-10)	10	10	10	9	10	10	10	10	10
Total Score	81	82	84	84	82	84	84	81	62
Percent Comparability to Reference Station	100	101	104	104	101	100	100	100	76

Assessment Category	Percent of Comparability
Comparable to Reference	≥ 90%
Supporting	75 -88%
Partially Supporting	60 -73%
Non-Supporting	≤ 58%

[1] Indicates Reference Station for that area
NOTE: EPA RBP Habitat Assessment not appropriate for pond stations.

Table A-9.2
RBP-I Macorbenthos Data

Taxon	TDTS STATION IDENTIFICATIONS								
	HMC-1	HMC-2	HMC-3	HMC-4	HMC-5	GRC-1	GRC-2	HHC-1	HHC-2
Oligochaeta									
Isopoda									
Amphipoda									
Decapoda	C	C	C	R				C	R
Insecta									
Anisoptera	C	C	C	R	C	C	C	C	A
Zygoptera	C		C						
Hemiptera									
Coleoptera	C		C		C	A	A	A	R
Diptera									
Tipulidae									
Tabanidae	R			C		R		C	
Culicidae						C			R
Chironomidae				C					A
Plecoptera	C		C	A	C	C	C		
Ephemeroptera	C	C							
Tricoptera									
Gastropoda			C						

Table A-9.3
RBP-III Habitat Assessment Data

Habitat Parameter (RANGE OF VALUES)	TDTS STATION IDENTIFICATIONS								
	HMC-1[1]	HMC-2	HMC-3	HMC-4	HMC-5	GRC-1[1]	GRC-2	HHC-1	HHC-2
Bottom Substrate (0-10)	10	16	11	14	15	16	15	12	6
Embeddedness (0-15)	2	2	2	0	2	5	2	2	2
Stream Flow and/or Stream Velocity (0-20)	5	11	16	18	16	16	16	4	5
Channel Aeration (0-15)	9	11	8	15	11	12	8	11	8
Bottom Scouring and Deposition (0-15)	12	12	9	12	11	12	10	12	7
Pool/Riffle or Run/Bend Ration (0-15)	8	12	11	12	13	11	10	13	8
Bank Stability (0-10)	9	9	9	8	10	9	8	9	8
Bank Vegetative Survey (0-10)	8	9	10	10	9	10	9	8	8
Streamside Cover (0-10)	5	9	8	8	8	8	8	8	7
Total Score	68	91	84	97	95	99	86	79	59
Percent Comparability to Reference Station	100	134	124	143	140	100	87	100	75

Assessment Category	Percent of Comparability
Comparable to Reference	≥ 90%
Supporting	75 -88%
Partially Supporting	60 -73%
Non-Supporting	≤ 58%

[1] Indicates Reference Station for that area
NOTE: EPA RBP Habitat Assessment not appropriate for pond stations.

Table A-9.4
RBP-III Benthos (Riffle/Run) Data

Class	Order	Family	Genus	Species	Station Identification																
					GaP-1	GaP-1a	GrC-1	GrC-2	HHC-1	HHC-2	HMC-1	HMC-2	HMC-3	HMC-4	HMC-4D	HMC-5	BeP-1	BeP-1a	BeP-2	HOP-1	REP-1
Crustacea	Amphipoda	Gammaridae	Gammarus						1	3								6	17	4	1
Crustacea	Decapoda	Cambaridae	Orconectes							1											
Crustacea	Decapoda						1	1	1			1						1			
Gastropoda		Physidae								1											
Crustacea	Isopoda																	2	58	1	
Crustacea	Isopoda	Asellidae	Caecidotea					1													
Hirudnea	Rhynchobdellidae	Glossiphomidaae	Helobdella	stagnalis				1		5				8	2	4		1	1	3	
Insecta	Coleoptera	Elmidae								1								1	1		
Insecta	Coleoptera	Hydrophilidae	Berosus																8		
Insecta	Coleoptera	Scirtidae															16	5	1		
Insecta	Diptera	Ceratopegonidae	Palopamoyia		1	3				1											
Insecta	Diptera	Chironomidae			8				2		2			4			210	8	52	12	9
Insecta	Diptera	Chironomidae	Polypedellum			1		1		5		2	1		2						1
Insecta	Diptera	Chironomidae	Tribelas		2	3										1			1		
Insecta	Diptera	Chironomidae	Chironomus			1							2				2	1	92	7	
Insecta	Diptera	Chironomidae	Xenochironomus										1	1							
Insecta	Diptera	Chironomidae	Endochironomus																	11	1
Insecta	Diptera	Chironomidae	Goeldichironomus		22																
Insecta	Diptera	Chironominae	Microspectra			10											2				
Insecta	Diptera	Chironomidae	Dicrotandpes														2				2
Insecta	Diptera	Chironomidae	Glyptepeudipes														7	11	2		
Insecta	Diptera	Chironomidae	Paratendipes														41				
Insecta	Diptera	Chironomidae	Microtendipes																3		
Insecta	Diptera	Chironomidae	Cryptochironomus																		1
Insecta	Diptera	Tipulidae									1			5			1				
Insecta	Diptera	Tauypodae	Procaldus							16											
Insecta	Diptera	Orthocladidae	Cricotopus							12											
Insecta	Diptera	Simulidae								11	2			1	3						
Insecta	Diptera	Tanypedidae										2									
Insecta	Diptera	Tanypedidae	Procladius		20	65								1			9	2	55	7	
Insecta	Diptera	Tanypedidae	Dialmeabatista			5															
Insecta	Diptera	Chilenydae																			
Insecta	Diptera	Tabanidae			1													2	2		
Insecta	Ephemoptera	Caenidae	Caenis														1				
Insecta	Ephemoptera	Leptophillidae	Brakptophibia							2											
Insecta	Ephemoptera	Leptophillidae	Paraleptophidae								2			1							
Insecta	Ephemoptera	Ephemerillidae	Ephemerella							13											
Insecta	Ephemoptera	Ephemerillidae	Dannella																43		
Insecta	Ephemoptera	Heptagenilidae	Stenenora				1	9		1	2		2	6		1					
Insecta	Ephemoptera	Baetidae											1								
Insecta	Ephemoptera	Baetidae	Centropilum					1		7				1					28		
Insecta	Ephemoptera											1							4		
Insecta	Hemiptera	Belestoniatidae	Belestoma																3		
Insecta	Hemiptera	Gelastocoridae	Nerthra																1		
Insecta	Hemiptera	Corixidae	Palmacorixa																4		
Insecta	Megaloptera	Corydalidae	Corydalus									1		1							
Insecta	Odonata	Aeshnidae	Boyeria								2										
Insecta	Odonata	Gomphidae	Gomphus					1	1			1									
Insecta	Odonata	Cordulidae	Tetragenenta					1										1	1		
Insecta	Odonata	Macromiidae	Macromia					1													
Insecta	Odonata	Libellulidae	Orthemis			3													3		
Insecta	Odonata	Libellulidae	Leucomhua															2	4		
Insecta	Odonata	Coenagrionidae	Argia																12		2
Insecta	Odonata	Lestidae	Lestes																17		

Table A-9.4
RBP-III Benthos (Riffle/Run) Data

Class	Order	Family	Genus	Species	Station Identification															
					GaP-1	GaP-1a	GrC-1	GrC-2	HHC-1	HHC-2	HMC-1	HMC-2	HMC-3	HMC-4	HMC-4D	HMC-5	BeP-1	BeP-1a	BeP-2	HOP-1	REP-1
Insecta	Plecoptera	Perlidae	Bebneuria											4							
Insecta	Plecoptera	Perlidae	Eccoptura											2	2						
Insecta	Plecoptera	Perlidae	Phasganephora					2				1									
Insecta	Plecoptera	Perlidae	Perinella												1						
Insecta	Plecoptera	Perlidae	Perlesta				1	1		8	2										
Insecta	Trichoptera										2				1						
Insecta	Trichoptera	Philopatamidae	Cymarra				1			2			1	3							
Insecta	Trichoptera	Hydropsychidae	Macrostenium												1						
Insecta	Trichoptera	Hydropsychidae	Hydropsyche						3			1									
Insecta	Trichoptera	Limnephilidae							1			1									
Insecta	Trichoptera	Limnephilidae	Pseudostenophylax						1												
Oligochaeta	Tubicifida	Tubicifidae				4			3	3							7	11	2	5	1
Tubellaria	Trichadidae	Planrlidae	Dugesia																		57
				Number of Organisms	54	95	4	20	13	92	15	11	8	38	12	6	298	54	415	50	75

Table A-9.5
RBP-III Benthos (CPOM) Data

Class	Order	Family	Genus	Species	Station Identification															
					GaP-1	GrC-1	GrC-2	HHC-1	HHC-2	HMC-1	HMC-2	HMC-3	HMC-4	HMC-4D	HMC-5	BeP-1	BeP-1a	BeP-2	HOP-1	REP-1
Crustacea	Amphipoda	Gammaridae	Gammarus						6	4						4	1	6	9	
Crustacea	Decapoda	Cambaridae	Orconeetes						1											
Crustacea	Cladocera															7	7	64		
Gastropoda		Ancylidae	Ferrissia						1											
Crustacea	Isopoda	Asellidae	Caeciodotea					1	1									11	1	
Insecta	Coloeptera	Elmidae				5	1		1		5				3	5	5	1	1	
Insecta	Coloeptera	Scirtidae																2		
Insecta	Coloeptera	Dysticidae			1															
Insecta	Diptera	Tipulidae			2	1	1		3	2		1	1							
Insecta	Diptera	Chironomidae	Tribelos		1				1				11	6						
Insecta	Diptera	Chironomidae			4					4	6			14	1	17	54	38	4	6
Insecta	Diptera	Tanypedidae	Procladius		3			2	12	4	1			5	1	28	19	6		9
Insecta	Diptera	Chironomidae	Cryptochironomus		1								1							
Insecta	Diptera	Chironomidae	Endochironomus		1		4				8								1	
Insecta	Diptera	Orthocldidae	Xylotopus			1							1							
Insecta	Diptera	Simulidae				1			1	1		1		1						
Insecta	Diptera	Chironomidae	Chironomus				1									5	19	145		3
Insecta	Diptera	Chironomidae	Polypedium						2				1							9
Insecta	Diptera	Tanytarsini	Microspectra						3			1								
Insecta	Diptera	Orthocladidae	Cricotopus							1				2						
Insecta	Diptera	Chironomidae	Xenochironomus							3										
Insecta	Diptera	Orthocladidae	Psectrocladus									1								
Insecta	Diptera	Ceratopogaridae	Palopmyia											2		7		1		
Insecta	Diptera	Chironomidae	Dicrotendipes													1	3	8		1
Insecta	Ephemoptera	Heptagridae	Stenonas			4			1	9	1	1			1					
Insecta	Ephemoptera	Baetidae	Centroptilium						7											
Insecta	Ephemoptera	Leptophlebidae	Paraleptophlebia							3								3		
Insecta	Ephemoptera	Ephemerellidae	Ephemerella													1		13	1	
Insecta	Ephemoptera	Baetidae														1				
Insecta	Ephemoptera						1													
Insecta	Hemiptera	Corixidae	Palmacorixa													1		1		
Insecta	Odonata												1							
Insecta	Odonata	Macromidae	Macromia			1														
Insecta	Odonata	Aeshidae	Boyeria			1														
Insecta	Odonata	Cordulidae	Tetragoneuria				1													
Insecta	Odonata	Gomphidae	Gomphus								1									
Insecta	Odonata	Gomphidae	Progamphs									1								
Insecta	Odonata	Libelluidae	Orthemis													4	2			
Insecta	Odonata	Libelluidae	Leucorrhina															5		
Insecta	Odonata	Coenagrienidaae														1				
Insecta	Odonata	Coenagrienidaae	Argia															4		
Insecta	Odonata	Lestidae	Lestes															2		
Insecta	Odonata	Lestidae																		1
Insecta	Plecoptera	Perlidae	Belomeria			1					1	1								
Insecta	Plecoptera	Perlidae	Perlesta						1	4		2			1					
Insecta	Trichoptera								1	3										
Insecta	Trichoptera	Limnephilidae				1			4											
Insecta	Trichoptera	Philopotanidae	Chamarra			1				1		1								
Insecta	Trichoptera	Beraridae	Beraea					9				1	3	1						

Table A-9.5
RBP-III Benthos (CPOM) Data

Class	Order	Family	Genus	Species	Station Identification															
					GaP-1	GrC-1	GrC-2	HHC-1	HHC-2	HMC-1	HMC-2	HMC-3	HMC-4	HMC-4D	HMC-5	BeP-1	BeP-1a	BeP-2	HOP-1	REP-1
Insecta	Trichoptera	Hydropsychidae	Hydropsyche									1		1						
Insecta	Trichoptera	Limnephilidae	Pseudostenephlax															1		
Tubellaria	Tricladidae	Planariidae	Dugesia																	29
Oligochaeta	Tubicifida	Tubicifidae			15			2					1	1	2	13	7			6
Oligochaeta	Lumbriculidae	Lumbriculis					1													
			Number of Organisms		28	17	10	14	46	39	23	12	20	33	9	95	117	311	17	64

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