GW -

# REPORTS

# YEAR(S):



#### STORMWATER POLLUTION PREVENTION PLAN

#### MILAGRO GAS CONDITIONING PLANT

Williams Field Services Company

August 2001

## **Table of Contents**

| 1.0 INTRODUCTION                                    | 1    |
|---|------|
| 1.1-Site Name and Location                          | 1    |
| 1.2 Site Operator                                   | 1    |
| 1.3 Plan Organization                               | 1    |
| 1.4 Industrial Activity Description                 | 1    |
| 1.5 Consistency with Other Plans                    | . 2  |
| 2.0 EXISTING CONDITIONS                             | 3    |
| 2.1 Facility Description                            | 3    |
| 2.2 Storm Water Drainage Patterns                   | 4    |
| 2.3 Non-Storm Water Discharges                      | 5    |
| 2.4 Prior Spills and Leaks                          | 6    |
| 2.5 Endangered Species                              | 6    |
| 2.6 Historic Properties                             | 7    |
| 3.0 STORM WATER MANAGEMENT CONTROLS                 | 8    |
| 3.1 Pollution Prevention Team                       | 8    |
| 3.1.1 SWP3 Coordinator                              | 8    |
| 3.1.2 SWP3 Manager/Administrator                    | 9    |
| 3.1.3 SWP3 Inspector                                | 10   |
| 3.2 Risk Identification and Assessment              | 10   |
| 3.3 Best Management Practices                       | 12   |
| 3.3.1 Good Housekeeping                             | 12   |
| 3.3.2 Preventive Maintenance                        | 12   |
| 3.3.3 Spill Prevention and Response Procedures      | . 13 |
| 3.3.4 Inspections and Evaluation                    | 13   |
| 3.3.5 Sediment and Erosion Control                  | 13   |
| 3.3.6 Process and Materials Management              | 14   |
| 3.3.7 Employee Training                             | 14   |
| 4.0 IMPLEMENTATION                                  | 15   |
| 4.1 Best Management Practice (BMP) Implementation   | 15   |
| 4.2 Visual Monitoring                               | 15   |
| 4.3 Employee Training                               | 16   |
| 4.4 Record Keeping & Internal Reporting Procedures  | 16   |
| 4.5 Annual Comprehensive Site Compliance Evaluation | 17   |

4.5 Annual Comprehensive Site Compliance Evaluation

i

#### List of Tables

3

 Table 1
 Product and Waste Storage Tanks

## **List of Figures**

Figure 1Site Vicinity/Topographic MapFigure 2Site Plan

Figure 3 Drainage Plan

1

## Appendices

Appendix ASWP3 DocumentationAppendix BInspection and Monitoring WorksheetsAppendix CStorm Water Visual Monitoring Procedure

#### **Storm Water Pollution Prevention Plan Certification**

Name of Facility: Milagro Gas Conditioning and Cogeneration Plant

**Type of Facility:** The Milagro Plant is a natural gas conditioning and cogeneration plant. The plant's Standard Industrial Code number is 1321. In the National Pollutant Elimination System (NPDES) Multi Sector General Permit (MSGP), the plant is categorized as a Sub-sector 2-Natural Gas Liquids Facility in Sector I-Oil and Gas Extraction and Refining.

**Location:** Milagro Gas Conditioning Plant is located in the W/2 of SE/4, Section 12, Township 29 North, Range 11 West, in San Juan County, New Mexico, in approximately 2.5 miles east of Bloomfield, New Mexico.

Site Address: 192 County Road 4900 Bloomfield, NM 87413

Storm Water Pollution Prevention Plan Certification: I certify under penalty of law that this document and all attachments were prepared under my direction or supervision in accordance with a system designed to assure that qualified personnel properly gathered and evaluated the information submitted. Based on my inquiry of the person or persons who manage the system, or those persons directly responsible for gathering the information, the information submitted is, to the best of my knowledge and belief, true, accurate, and complete. I am aware that there are significant penalties for submitting false information, including the possibility of fine and imprisonment for knowing violations.

Name: JEFFREY D. BAUMER

Signature:

Date: 8/15/01

#### 1.0 INTRODUCTION

This Storm Water Pollution Prevention Plan (SWP3) is developed for implementation at the Milagro Gas Conditioning and Cogeneration Plant for storm water discharges associated with industrial activity. This SWP3 shall be implemented to fulfill the requirements of Section 405 of the Water Quality Act of 1987, which added Section 402(p) to the Clean Water Act. This section dictated that the Environmental Protection Agency (EPA) establish regulations setting forth National Pollutant Discharge Elimination System (NPDES) permit application requirements for storm water discharges associated with industrial activity and discharges from municipal separate storm sewer systems. The SWP3 outlined herein is designed to fulfill permit requirements.

1.1 Site Name and Location

Milagro Gas Conditioning and Cogeneration Plant is located in the W/2 of SE/4, Section 12, Township 29 North, Range 11 West, in San Juan County, New Mexico, approximately 2.5 mile east of Bloomfield, New Mexico. A site location map is attached (USGS 7.5 Min. Quadrangle: Bloomfield, New Mexico) as Figure 1. A facility plot plan is attached as Figure 2.

The site address is: 192 County Road 4900 Bloomfield, NM 87413

1.2 Site Operator

The Milagro Gas Conditioning Plant is operated by Williams Field Services (WFS).

1.3 Plan Organization

This SWP3 is consistent with the Federal Water Pollution Control Act and the New Mexico 401 Water Quality Certification. Storm water discharges are allowed under the general National Pollutant Discharge Elimination System (NPDES) Storm Water Multi-Sector General Permit (MSGP). On March 28, 2001, the EPA issued permit number NMR05A883 for the facility. The active date of permit coverage was March 14, 2001.

This plan concentrates on the identification of personnel responsible for implementation of the SWP3, identification of potential pollutants, description of structural controls to prevent pollution of storm water pollution, maintenance, inspection and evaluation, and record keeping.

1.4 Industrial Activity Description

The Milagro Plant is a natural gas conditioning and cogeneration plant. The plant's Standard Industrial Code number is 1321. In the MSGP, the plant is categorized as a Sub-sector 2-Natural Gas Liquids Facility in Sector I-Oil and Gas Extraction and Refining.

#### 1.5 Consistency with Other Plans

This SWP3 has been developed in conjunction with the existing environ-mental and operating plans listed below.

- The Milagro Gas Conditioning Plant currently implements a Spill Prevention Control and Countermeasure (SPCC) Plan, 42.13.001. This SPCC Plan is in accordance with Section 311 of the CWA, CFR Title 40, Part 112.7, State and Local Government Requirements.
- Emergency Plan for Milagro (ref. Emergency Operating Procedures, 42.01.001 and 42.01.002).
- Discharges or Spills of Oil or Hazardous Substances (ref. Operating and Maintenance Procedure, 21.10.020).
- New Mexico Oil Conservation Division Discharge Plan

#### 2.0 EXISTING CONDITIONS

#### 2.1 Facility Description

This facility, built in 1991, is a natural gas conditioning and cogeneration plant. The conditioning plant is designed to remove carbon dioxide and water from raw natural gas. Plant processes include gas dehydration using triethylene glycol,  $CO_2$  removal by contacting natural gas with methyldiethanolamine (MDEA) and glycol and MDEA regeneration. The processes used to separate hydrocarbons from field gas includes compressors, heat exchangers, separators, dehydrators, power generating equipment, aboveground storage tanks and other supporting equipment.

The cogeneration plant generates electricity using natural gas fuel. The exhaust gas heat is used to generate steam for use in the gas conditioning plant. In addition, there are various storage tanks, support structures and ancillary equipment. The product and waste storage tank descriptions, capacities, and locations are summarized in Table 1 and Figure 2.

| #          | Number<br>of Tanks | Capacity<br>(approximate) | Description                         | Drainage Basin |
|------------|--------------------|---------------------------|-------------------------------------|----------------|
| [          |                    |                           |                                     |                |
| <u>(a)</u> | 1                  | 1,175 gallon              |                                     | Area 2         |
| <u>(b)</u> | 1                  |                           | transformer                         | Area 2         |
| (c)        | 1                  |                           | ambitrol/water tank                 | Area 2         |
| (d)        | 1                  |                           | triethylene glycol tank             | Area 2         |
| (e)        | 2                  | 250 barrel                | amine tanks                         | Area 2         |
| (f)        | 1                  | 500 barrel                | amine/water tank                    | Area 2         |
| (g)        | 1                  | 100 barrel                | amine/water tank                    | Area 2         |
| (h)        | 1                  | 300 gallon                | solvent tank                        | Area 2         |
| (i)        | 1                  | 400 gallon                | polymer tank                        | Area 2         |
| [ (j)      | 1                  | 400 gallon                | cortrol tank                        | Area 2         |
| (k)        | 1                  | 400 gallon                | steamate tank                       | Area 2         |
| (1)        | 2                  | 100 barrel                | triethylene glycol /<br>water tanks | Área 2         |
| (m)        | 1                  | 500 gallon                | gasoline/diesel tank                | Area 2         |
| (n)        | 2                  | 456 gallon                | transformers                        | Area 3a        |
| (0)        | 4                  | 233 gallon                | transformers                        | Area 3a        |
| (p)        | 2                  | 2,300 gallon              | lube oil tanks                      | Area 3a        |
| (q)        | 1                  | 1,250 gallon              |                                     | Area 3a        |
| (r)        | 2                  | 4,233 gallon              | transformers                        | Area 3b        |
| (s)        | 3                  |                           | evaporation tanks                   | Area 8         |
| (t)        | 1                  | 100 barrel                | used oil tank                       | Area 8         |
| (u)        | 1                  | 250 barrel                | slop oil tank                       | Area 8         |
| (v)        | 1                  | 1,000 gallon              | used oil tank                       | Area 8         |
| (x)        | 1                  | 250 barrel                | inlet/outlet filter<br>liquids tank | Area 8         |

# TABLE 1PRODUCT AND WASTE STORAGE TANKS

#### 2.2 Storm Water Drainage Patterns

The Milagro Plant occupies approximately 40 acres of land with the elevation of the facility ranging from approximately 5,720 to 5,645 feet above mean sea level. Regional surface water drainage is to the south-southeast along Hare Canyon Wash towards San Juan River. Milagro Plant storm water drainage consists of Hare Canyon Wash located west of the facility and an unnamed arroyo located south of the facility. The unnamed arroyo drains into Hare Canyon Wash approximately 1 mile south of Milagro Plant.

Nine discrete storm water drainage basins were identified at the facility during the site inspection. Storm water drainage basins. Figure 3 identifies each of the site drainage basins and their points of storm water discharge. Each area is discussed in the following sections:

• Plant Process Area Gas Tranes 1 through 4 (Area 1a)

This process area includes Gas Tranes 1 through 4. Storm water flows towards the west. In the Trane #1 area, a surface drain system collects storm water. A french drain system collects storm water in the area of Tranes 2 through 4. The surface drain and french drain systems discharge southwest of the area.

• Plant Process Area Gas Trane 5 (Area 1b)

This process area includes Gas Trane 5. Storm water flows towards the south into a drainage conveyance. The drainage conveyance discharges to the south portion of the property.

• Utility Area (Area 2)

The area consists of glycol regeneration skids, water treating building, auxiliary boiler, generator building, air compressor building, and storage tank area. Storm water from this area flows northwest to a conveyance consisting of piping and ditch. The conveyance discharges west of the facility into the Hare Canyon Wash.

• Cogeneration Area (Area 3a)

The area consists of electric generators powered by a natural gas fired turbine. Storm water from this area flows west to diversion conveyance consisting of piping and ditch. The conveyance discharges west of the facility into the Hare Canyon Wash.

• Switch Yard (Area 3b)

The area consists of electric switchgear equipment. Storm water from this area flows west to conveyance consisting of a wide ditch. The conveyance discharges west of the facility into the Hare Canyon Wash.

• Wastewater Evaporation Pond Area (Area 4)

The evaporation pond area is located at the immediately west of the main facility. The area consists of three evaporation ponds. The evaporation ponds capture storm water runoff in this area. No storm water is discharged from this area.

• Slope Drainage Area (Area 5)

Storm water from the western slope collects in a ditch and is directed either west towards Hare Canyon Wash or into the south discharge area.

• South Embankment Area (Area 6)

Storm water is collected by a conveyance along the lip of the embankment to a drainage conveyance that discharges at the base of the embankment.

• Plant Parking Area (Area 7)

Storm water flows from this area towards the southwest and enters a conveyance towards the west. Storm water is discharged into an unnamed ditch that flows into Hare Canyon Wash.

• North 40 Lay-Down Area (Area 8)

Storm water from this area flows west to Hare Canyon Wash.

• South Discharge Area (Area 9)

This area receives storm water from Areas 1a, 1b, 5, 6 and 7. Storm water flows southwest and discharges through gate valve and conduit.

2.3 Non-Storm Water Discharges

The SWP3 team conducted a visual assessment of each of the storm water outfalls that discharge off-site on June 26, 2001. Storm water outfalls were inspected for evidence of non-authorized discharges (staining, odors, discharge pipes, etc.) from the facility. There were no visual indications of; non-authorized discharges observed. Certification of the assessment for the presence of non-authorized discharges is presented in Appendix A.

Non-storm water discharges that potentially flow to the storm water management system include the following:

- Discharges from emergency fire-fighting activities, fire equipment flushings and tests,
- Air conditioning condensate,
- Routine exterior building wash downs (non-detergent),
- Potable water from fire water pumps, and
- Routine pavement wash downs (non-detergent) where spills or leaks have not occurred.

The discharges identified above are authorized under the MSGP based on their identification in this SWP3. Non-storm water discharges that are not authorized under the MSGP include the following:

- Cooling tower water,
- Process area waste water,
- Compressor wash water,
- Maintenance slab wash water, and
- Sanitary wastewater.

With the exception of sanitary wastewater that flows to the facility septic system, the above referenced, non-authorized discharges flow to the evaporation ponds via underground piping. Wastewater flows from the process area floor drains flow to an oil/water separator prior to discharging to the evaporation ponds.

2.4 Prior Spills and Leaks

Releases at the Milagro Plant are recorded and documented on Spill Reports maintained in the facility's environmental files. A reportable quantity (RQ spill occurred at the facility in late January 1999 that triggered the requirement for this SWP3. A summary of this spill is summarized below. No other significant spill events in the past three years.

• In late January 1999, water was observed flowing from a junction box approximately 300 feet west of the Milagro Plant Office Building. The estimated volume of discharged water was 9,540 gallons. The discharged water consisted of a mixture of septic and wastewater. Significant settlement of the building's foundation caused failure of the drain lines. The drain line were tested and repaired. The affected soils were assessed and clean up was performed.

#### 2.5 Endangered Species

The NOI requires MSGP applicants to identify whether threatened and endangered (T&E) species are present in proximity to storm water discharges. The threatened and endangered species have been identified in San Juan County are the Bald Eagle, Mexican-Spotted Owl, the Back-Footed Ferret and Knowlton Cactus. According to the Threatened and Endangered Species Conflict Map located at the Bureau of Land Management (BLM) Farmington, New Mexico office, no threatened or endangered species were identified in the Milagro Plant area.

Based on existing storm water discharges under the MSGP and the established storm water management system, only minimal land disturbing activities (i.e. berm repairs, ditch cleanings) are anticipated at this time. With the exception of the incidental occurrence of Bald Eagles, no T&E species are known to be "In Proximity" to the facility. Based on visual inspection, the industrial nature of the Milagro Plant, and minimal disturbance of natural habitat, adverse effects to these species are not likely.

#### 2.6 Historic Properties

The NOI requires MSGP applicants to identify historical properties that are present in proximity to storm water discharges. The historical property must be either listed or eligible for listing on the National Register of Historical Places. A search of the National Register Information System did not reveal any historical properties in the proximity of Milagro Plant. A subsequent visual inspection of the facility did not identify any historic properties that would be affected by implementation of this plan.

#### 3.0 STORM WATER MANAGEMENT CONTROLS

#### 3.1 Pollution Prevention Team

The Pollution Prevention Team consists of specific individuals responsible for the development and revision, implementation, and maintenance of the SWP3. The Team will consist of management personnel responsible for overseeing the plan, conducting quarterly inspections, conducting annual evaluations, and preparing reports. The Team will also include on-site technical personnel responsible for the day-to-day activities and maintenance of the SWP3. Team members are comprised of the following:

| Name                           | Team Position      | Telephone No.                    |
|--------------------------------|--------------------|----------------------------------|
| Mark Bareta                    | Coordinator        | (505) 632-4634                   |
| Eric Edmanson<br>Bryan Salazar | Manager/Administer | (505) 632-4613<br>(505) 632-4735 |
| J.D. Brow or<br>Designee       | Inspector          | (505) 632-4607                   |

#### 3.1.1 SWP3 Coordinator

The SWP3 Coordinator for the Milagro Gas Processing Plant will be located at the San Juan Area office of Williams Field Services (WFS) in Bloomfield, New Mexico. The Coordinator will be responsible for the following SWP3 activities:

- <u>Coordination with Regulatory Agencies</u>: The Coordinator will be prime contact with the regulatory agencies involved in the NPDES permit process and shall be responsible for compiling and submitting requested information.
- <u>Maintenance of Required Documentation</u>: The Coordinator will be responsible for ensuring that required documentation is maintained at the head office and on-site, including records of inspection, maintenance, spills/leaks, annual evaluation for compliance, revisions to the plan, and activities of Team members, records of training, and monitoring records.
- <u>Review of Inspection and Maintenance Schedule</u>: The Coordinator will review the inspection and maintenance schedule at least once annually to ensure that it fulfills the purpose of the plan. The schedule will be discussed with the Manager/Administrator, and any revisions agreed upon will be documented. The Coordinator shall also review inspection and maintenance reports to ensure that the schedule is being implemented.
- <u>Review of Employee Training Program</u>: The Coordinator shall review the schedule and content of the Employee training program to ensure that it fulfills the intent of the Plan and is keeping with the latest regulations and practices. The Coordinator will be available to conduct training efforts and assist the Manager/Administrator in formulation of training programs.
- <u>Annual comprehensive Site Compliance Evaluation</u>: The Coordinator shall conduct an annual on-site compliance evaluation with the Manager/Administrator. The Coordinator

shall be responsible for intermediate evaluations if observations during the annual evaluation indicate that the intent of the plan is not being fulfilled. The Coordinator shall prepare a report following the evaluation in accordance with the plan.

• <u>Revisions to the SWP3</u>: The Coordinator shall review the SWP3 on an annual basis and implement revisions or additions as needed to bring the plan current with the latest regulations and procedures and with any change in site conditions. These recommendations will be coordinated with the Manager/Administrator for implementation.

#### 3.1.2 SWP3 Manager/Administrator

The SWP3 Manager/Administrator shall visit the facility on a regular basis and shall report directly to the Coordinator. The Manager/Administrator shall be responsible for the following Plan activities:

- <u>Inspection and maintenance schedule</u>: The Manager/Administrator will be responsible for developing and implementing the inspection and maintenance schedule and for assigning the tasks related to the schedule to the Inspector. The Manager/Administrator shall review the inspection and maintenance reports prepared by the Inspector and approve the reports prior to submitting copies to the Coordinator. The Manager/Administrator will be responsible for ensuring that the schedule is implemented and proper documentation is completed. The Manager/Administrator will maintain a file of all documentation on-site. The Manager/Administrator shall ensure that any deficiency found during inspection is rectified in a timely manner and documented in the file.
- <u>Employee Training Program</u>: The Manager/Administrator shall be responsible for developing and implementing an employee-training program for site personnel involved in activities related to the plan. The Manager/Administrator shall ensure that all personnel are aware of permit and regulatory requirements associated with their activities and are familiar with the contents of the plan.
- <u>Annual comprehensive Site Compliance Evaluation</u>: The Manager/Administrator shall accompany the Coordinator during site evaluations to provide information and ensure that any deficiency found during evaluation is rectified in a timely and well-documented manner. The Manager/Administrator will maintain a record of evaluation reports and resulting actions on-site.
- <u>Spills and Leaks</u>: The Manager/Administrator shall respond immediately to spills and leaks and take appropriate action to repair and mitigate such events. The Manager/Administrator shall notify the Coordinator immediately when any event threatens to pollute the storm water system. The Manager/Administrator will be responsible for documenting all events, actions taken, and the extent, if any, of storm water contamination. This documentation shall be maintained on-site and a copy forwarded to the Coordinator.
- <u>Revisions to the SWP3</u>: The Manager/Administrator shall review the SWP3 on an annual basis and recommend revisions or additions as needed to bring the plan current with the latest regulations and procedures and with any change in site conditions. These recommendations will be coordinated with the Coordinator for implementation.

#### 3.1.3 SWP3 Inspector

The SWP3 Inspector shall be located on-site on a day-to-day basis and shall report directly to the Manager/Administrator. The Inspector shall be responsible for the following SWP3 activities:

- <u>Inspection and Maintenance</u>: The Inspector shall conduct routine scheduled inspections and prepare a report of the inspection to be submitted to the Coordinator. The Inspector shall oversee routine scheduled maintenance and request any replacement, repairs, or unscheduled maintenance needed and report to the Manager/Administrator on all work performed. The inspector shall perform storm water monitoring according to the requirements of this plan.
- <u>Employee Training Program</u>: The Inspector shall participate in the employee-training program and shall alert site personnel to any activity not in conformance with the intent of the plan. Any activity not in conformance with the plan shall be reported to the Manager/Administrator.
- <u>Spills and Leaks</u>: The Inspector shall immediately report any spill or leak according to the WFS SPCC Plan and Emergency Operating Procedures and take appropriate action to remedy the event, contain any spilled substance, and mitigate any damage. The Inspector shall provide a written report to the Manager/Administrator following any spills or leaks, describing the event, personnel involved, cause, and actions taken.
- <u>Revisions to the Plan</u>: The Inspector shall be thoroughly familiar with the requirements of the SWP3 and may offer comments or recommendations to the Manager/Administrator concerning revisions or additions to the plan to better suit the site-specific conditions.
- 3.2 Risk Identification and Assessment

Eight drainage areas of the site, shown in Figure 3, have been assessed for their potential as storm water pollution sources. Preventative measures that have been implemented for each area are described below.

• Plant Process Area Gas Tranes 1 through 4 (Area 1a)

The potential pollutants are methyldiethanolamine (MDEA), glycol, and steam condensate. Concrete containment structures reduce the risk of amine releases to the ground surface.

• Plant Process Area Gas Trane 5 (Area 1b)

The potential pollutants are MDEA, glycol, and steam condensate. Concrete containment structures are located to reduce the risk of amine releases to the ground surface.

• Utility Area (Area 2)

The potential pollutants are MDEA, ambitrol, cortrol, dielectric oil, diesel, gasoline, polymer, solvent, steamate, and triethylene glycol. Product and waste storage tanks are located in this area. Potential leaks and spills associated with transfer activities from

these tanks could contribute pollutants to the storm water management system. The storage tanks have secondary containment capable of containing releases and preventing discharge of contents to the storm watercourses. Where experience indicates a reasonable potential for failure such as a tank overflow, rupture, or leakage, appropriate containment, diversionary structures, or equipment to prevent a discharge from reaching a watercourse are provided. In event of a spill, the SPCC Plan will be implemented.

• Cogeneration Area (Area 3a)

The potential pollutants are ambitrol, oil, glycol and steam condensate. The storage tanks have secondary containment capable of containing releases and preventing discharge of contents to the storm water system. In event of a spill, the SPCC Plan will be implemented.

• Switch Yard (Area 3b)

The potential pollutant is dielectric oil contained in the transformer. In event of a spill, the SPCC Plan will be implemented.

• Wastewater Evaporation Pond Area (Area 4)

The potential pollutants are oil and process wastewater. Process wastewater may consist of oil, glycol, ambitrol and amine. Since the tanks do not discharge to the surface, they are not a storm water pollution source. Storm water that falls or drains into the pond is evaporated along with the process wastewater. The evaporation tank levels are inspected daily. The storage tanks have secondary containment capable of containing releases and preventing discharge of contents to the storm water system. In event of a spill, the SPCC Plan will be implemented.

• Slope Drainage Area (Area 5)

The potential pollutants are oil, glycol and amine. A filter draining operation consisting of filter draining apparatus in a concrete containment and disposal dumpster could contribute petroleum hydrocarbons and process fluids. The filter draining operation is located with in a concrete-curbed area to contain runoff.

• South Embankment Area (Area 6)

The potential pollutant is steam condensate.

• Plant Parking Area (Area 7)

The potential pollutants are oil, glycol and septic wastewater from vehicle parking lot. The potential impact to storm water is minimal.

• North 40 Lay-Down Area (Area 8)

The potential pollutants are oil and process wastewater. Process wastewater may consist of glycol and amine. Equipment and material storage in this area could contribute petroleum hydrocarbons to off site drainages. Good housekeeping measures and enforcement of appropriate storage practices that include draining of equipment fluids prior to storage will minimize storm water impact

• Southwest Discharge Area (Area 9)

No potential pollutant sources are located in this area. In the event of a spill, the gate located in the southwest corner of the site will be closed to contain potentially impacted storm water.

#### 3.3 Best Management Practices

This section identifies the Best Management Practices (BMPs) currently being implemented to prevent or mitigate pollution to storm water.

#### 3.3.1 Good Housekeeping

All areas on-site are routinely maintained and kept free from waste material and debris. All waste material and debris is properly collected in appropriate container and disposed of on an on-going basis. Any temporary storage of potential pollutant sources is located in a contained hazardous material area with appropriate controls for storm water protection. All raw materials are protected from precipitation by storage in containers or coverage. All containers are clearly labeled with an accurate description of the substance contained. Personnel are to ensure that any reused containers are clearly marked with the current contents description and date. The previous descriptions shall be removed or voided. All containers are sealed; no open storage of liquid wastes is permitted.

#### 3.3.2 Preventive Maintenance

Periodic inspection and maintenance of storm water management devices as well as inspecting and testing facility equipment and systems to uncover conditions that could cause breakdowns or failures resulting in discharges of pollutants to surface waters, and ensuring appropriate maintenance of such equipment and systems.

- Open Channels: Periodic inspection of the open channels shall be performed at least once quarterly and after any large precipitation event. The channels shall be cleared of debris and sediment deposits and repaired and maintained as necessary.
- Culverts: Periodic inspection of the culverts shall be performed at least once quarterly and after any large precipitation event. Sediment deposits and debris shall be removed from the culverts, any erosion damage at entrance or outfall shall be repaired and extra protection considered if this is an on-going problem. Any damaged culvert shall be repaired or replaced.
- Containment Structures: Periodic inspection of the containment structures around process equipment, product and waste storage tanks, and the truck loading equipment shall be performed at least once quarterly and after any large precipitation event.

Process Wastewater System: Periodic inspection of the process wastewater system shall be performed at least once quarterly. The system will be evaluated and necessary maintenance

performed to ensure proper flow from the process area drains through the oil/water separator to the evaporation ponds.

#### 3.3.3 Spill Prevention and Response Procedures

Detailed policy and procedures for preventing, controlling, and reporting discharges or spills can be found in WFS Operating and Maintenance procedure, entitled Discharges or Spills of Oil or Hazardous Substances: Preventing, Controlling and Reporting of, 21.10.020 and the Milagro Plant SPCC Plan, 42.13.001.

#### 3.3.4 Inspections and Evaluation

In addition to or as part of the Annual Comprehensive Compliance Evaluation Report required by the permit, the qualified Inspector will conduct visual inspections of the Plant annually. The Inspector shall include quarterly visual inspections as part of his routine preventive maintenance site check. Records of annual inspections shall be maintained for a minimum of five years.

The objective of the visual inspections is to survey equipment and storage areas on a regular basis and to identify problems and ensure that appropriate actions are taken. A work order will be prepared describing corrective measures shall be used to ensure that appropriate actions are taken in response to the inspection.

Specifically, the Inspector will visually check for:

- Corroded or damaged drums, tanks, or pipes;
- Broken or breached earthen dikes;
- Broken or cracked concrete containment walls;
- Clogged catchment ponds;
- Collapsed or clogged culverts; and
- Stained soil.

#### 3.3.5 Sediment and Erosion Control

The following BMPs for sediment and erosion control are maintained and upgraded as needed to minimize the entrainment of fine-grained soils into storm water discharges.

- Maintain vegetative cover to the extent possible in exposed soil areas.
- Minimize soil erosion by reducing storm water runoff velocity and providing drainage pathways.
- Minimize exposure of bare soil areas to precipitation following any new construction or other ground disturbing activities. This can be accomplished by slope protection, flow diversions, and/or soil stabilization (mulching, matting, geotextiles).
- Allow sediments to settle out of storm water by utilizing the storm water detention pond.
- Ground stabilization at non-paved surfaces, such as parking, loading, unloading, and open areas. This may include re-grading the unpaved surface to remove ruts and erosion scars that have formed, followed by the application of gravel over non-paved areas.
- The majority of the plant process area has been covered with a layer of gravel to minimize erosion and potential contamination problems.
- Stabilize steep slopes.
- Minimize offsite runoff flows on to the facility by maintaining berms and diversions.

#### 3.3.6 Process and Materials Management

Materials management practices include the identification of toxic and hazardous substances stored on-site and the organization and handling procedures for these substances. By minimizing the exposure of materials to storm water, facilities can eliminate the possibility of storm water discharges coming into contact with pollutants. Specific practices emphasized include:

- Neatly organizing drums, boxes, and other containers for storage at designed storage areas;
- Identification of all substances stored and handled on-site;
- Development of handling procedures for these materials; and
- Revision of the SPCC and this Plan to include new materials stored or handled on-site.

#### 3.3.7 Employee Training

The SWP3 manager and inspector will conduct and coordinate employee training annually to inform and educate plant personnel about the SWP3. The following subjects will be covered:

- SPCC plan and associated spill response procedures;
- Good house keeping and preventative maintenance procedures;
- Proper tank filling, emptying, and material loading and unloading procedures;
- Drum handling, labeling and disposal procedures;
- Waste handling and disposal procedures; and
- Storm water monitoring and sampling procedures.

#### 4.0 IMPLEMENTATION

#### 4.1 Best Management Practices (BMP) Implementation

The Best Management Practices described in this plan have been implemented or will be initiated. These BMPs will be incorporated into the existing preventive maintenance schedule. Implementation will be monitored through inspections. The BMP, inspection interval, and responsible party are listed below.

| Best Management Practice                 | Inspection Interval | Responsible Party     |
|--|---------------------|-----------------------|
| Good Housekeeping                        | Annual              | Inspector             |
| Preventative Maintenance                 | Annual              | Inspector             |
| Spill Prevention and Response Procedures | Annual              | Inspector             |
| Visual Inspection                        | Annual              | Inspector             |
| Sediment and Erosion Control             | Annual              | Inspector             |
| Process and Materials Management         | Annual              | Inspector             |
| Employee Training                        | Annual              | Manager/Administrator |

#### 4.2 Visual Monitoring

Visual monitoring of storm water from each outfall are requirements of the MSGP. Visual monitoring of storm water discharge from the outfalls described in the section below must be conducted at least once per quarter to inspect storm water quality associated with snowmelt, or storm water runoff. Monitoring will be conducted during the following periods:

- 1<sup>st</sup> Quarter January through March,
- 2<sup>nd</sup> Quarter April through June,
- 3<sup>rd</sup> Quarter July through September, and
- 4<sup>th</sup> Quarter October through December.

Outfall sampling locations are shown in Figure 3 and include the following:

- Outfall 1 (Areas 1a, 1b, 4, 5, 6 and 9) Outfall location is the culvert located at the southwest boundary when flowing,
- Outfall 2 (Area 7) Outfall location to unnamed drainage when flowing,
- Outfall 3 (Areas 2, 3a, and 7) Outfall location to parking pavement when flowing, and,
- Outfall 4 (Areas 2, 3a, 3b, and 6) Outfall location to unnamed drainage when flowing.
- Outfall 5(Area 8) Outfall location to unnamed drainage when flowing.

Visual examinations are conducted during daylight hours within the first 30 minutes, or as soon as possible thereafter, after runoff or snowmelt begins discharging. All samples will be collected from the discharge resulting from a storm event that is greater than 0.1 inches in magnitude and that occurs at least 72 hours from the previously measurable (greater than 0.1 inches rainfall) storm event. The 72-hour storm interval is waived when the preceding measurable storm did not yield a measurable discharge, or it can be documented that less than 72-hour interval is representative for local storm events during the sampling period.

The grab sample must be collected in the first 30 minutes of the discharge. If the collection of a grab sample during the first 30 minutes is not possible, a grab sample can be collected during the first hour of the discharge, and the discharger must submit with the monitoring report a description of why a grab sample during the first 30 minutes was impracticable. A grab sample of the discharge is collected in a clear container. The visual examination is conducted in a well lit area during daylight hours and include any observations related to color, odor, clarity, floating solids, settled solids, suspended solids, foam, oil sheen, or other indicators of storm water pollution.

In some of the above-described areas, there are several outfalls within the same area that discharge offsite. A single sample for visual monitoring from one of the outfalls is deemed as a "representative discharge" based on the similar nature of industrial activities, type of materials, and management practices conducted within these areas.

Visual monitoring worksheets must be maintained onsite and filed with the SWP3. Blank worksheets are included in Appendix B. Information on the worksheets includes examination dates, times, personnel, nature of discharge (snowmelt, rainstorm, etc.), visual quality of the discharge, and probable sources of any observed pollution. If no discharges occur within a quarter then visual monitoring is not required and "no discharge" is recorded on the monitoring worksheets.

#### 4.3 Employee Training

An employee-training program shall be developed and implemented to inform personnel of storm water BMPs identified in this Plan. Training will be provided annually and records of said training shall be retained for three years. The employee training programs shall inform personnel responsible for storm water management and employees at all levels of responsibility of the components and goals of the storm water pollution prevention plan. Training addresses topics such as spill response, good housekeeping and material management practices.

The purpose of the training program is to inform and teach on-site personnel with regard to the components of the discharge regulations, permit, and storm water pollution prevention plan. The program will prepare individuals to effectively minimize and/or eliminate pollutants from entering the storm drainage system.

The goal of the program is to produce trained personnel who have the know-how to prevent spills and to respond safely and effectively, and who recognize and report potential storm water contamination situations.

#### 4.4 Record Keeping & Internal Reporting Procedures

All records of inspections and maintenance activities must be incorporated in Appendix C of this plan and maintained on-site. Documenting all inspections, training programs and incidents is an excellent preventive maintenance technique. A log will be developed and followed to record all maintenance, monitoring, and inspection activities with regard to the storm water drainage system. This log will become incorporated into the preventive maintenance log. Records of spills, leaks, or other pollutant discharges, inspections, monitoring events and maintenance activities must be retained for at least five years.

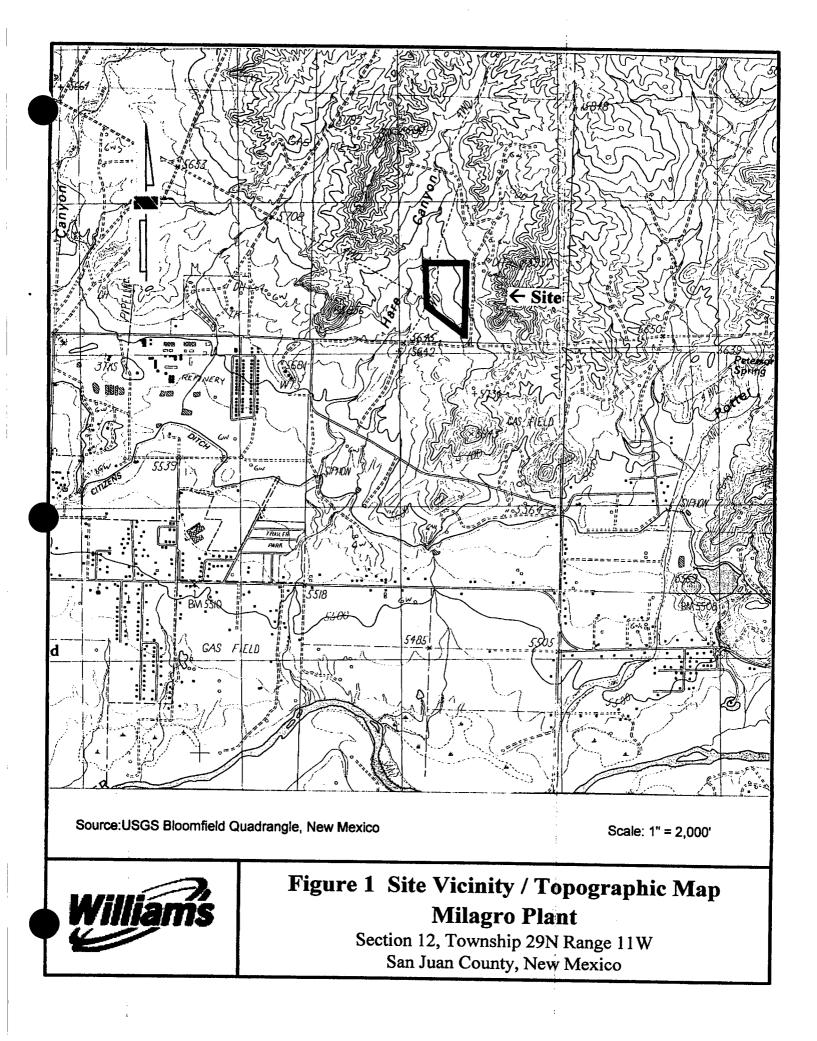
#### 4.5 Annual Comprehensive Storm Water Compliance Evaluation

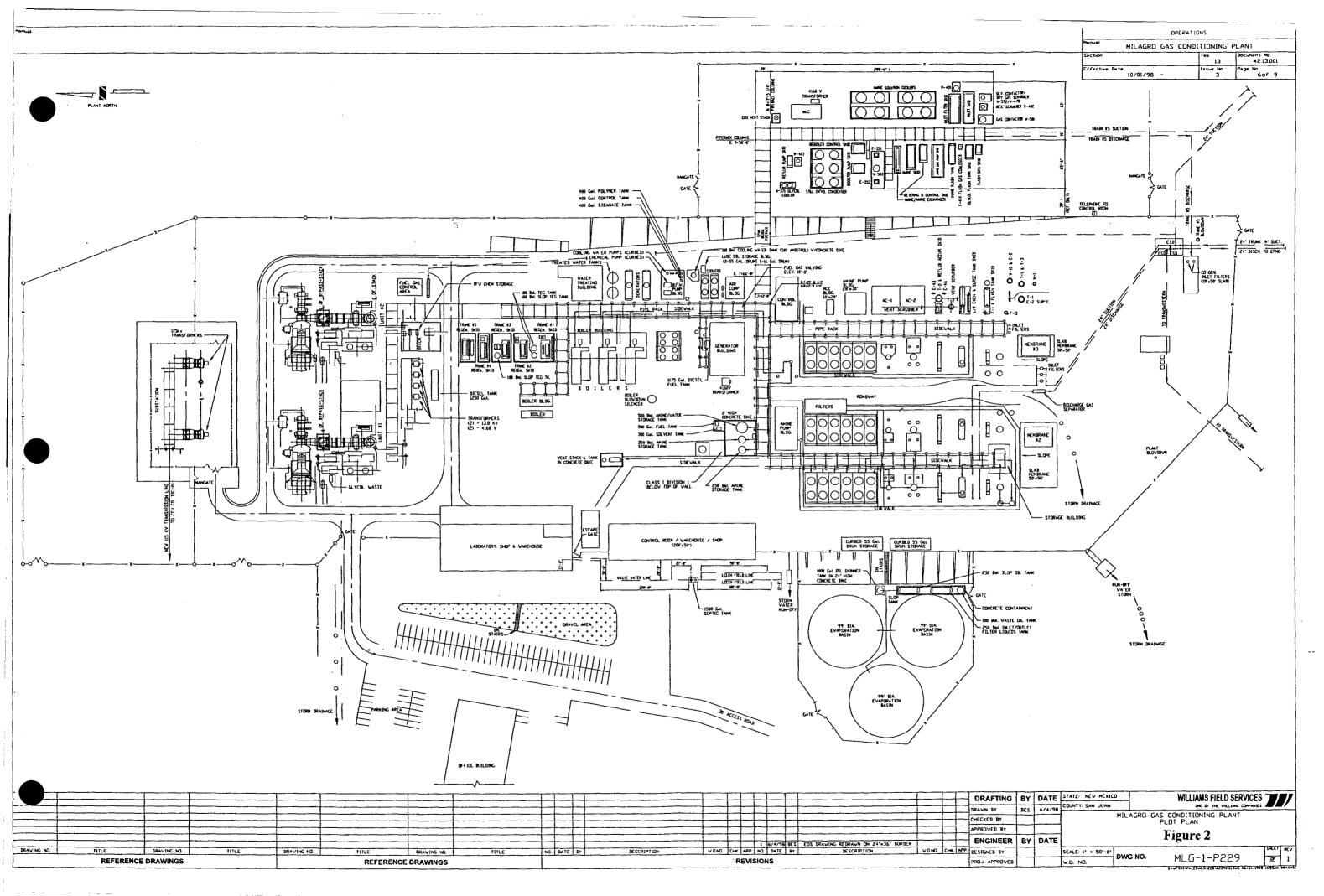
The Plan Coordinator and Plan Inspector shall conduct annual site compliance inspection. This annual inspection will provide a basis for evaluating the effectiveness of the SWP3. The inspection will evaluate measures to reduce pollutant loadings: structural storm water management measures, sediment and erosion control measures, and other structural pollution prevention measures identified in the SWP3. Measures shall be evaluated to determine whether they are adequate and properly implemented in accordance with terms of the permit or whether additional control measures are needed. A visual inspection of equipment needed to implement the SWP3, such as spill response equipment, shall be made. Areas contributing to any storm water discharge shall be visually inspected for evidence of, or the potential for, pollutants entering the drainage system. Specifically, the site evaluation will include:

- Inspect storm water drainage areas;
- Observe structural measures, culverts, earthen berms, and concrete dikes for proper operation;
- Evaluate effectiveness of storm water pollution prevention measures and BMPs;
- Review and evaluate quarterly monitoring data;
- Revise SWP3 to reflect new construction areas and changes in the storm water drainage system;
- Implement changes as required; and
- Complete and sign the Storm Water Compliance Evaluation Report.

A Storm Water Compliance Evaluation Report summarizing the scope of the inspection, personnel making the inspection, the date(s) of the inspection, major observations relating to the implementation of the SWP3, and actions taken shall be prepared. If the report describes deficiencies in pollution control structures or procedures, such deficiencies shall be corrected and the SWP3 shall be modified to reflect the required changes. The report shall be signed by the Coordinator and Inspector. The report shall be retained as part of the SWP3 for at least five years.

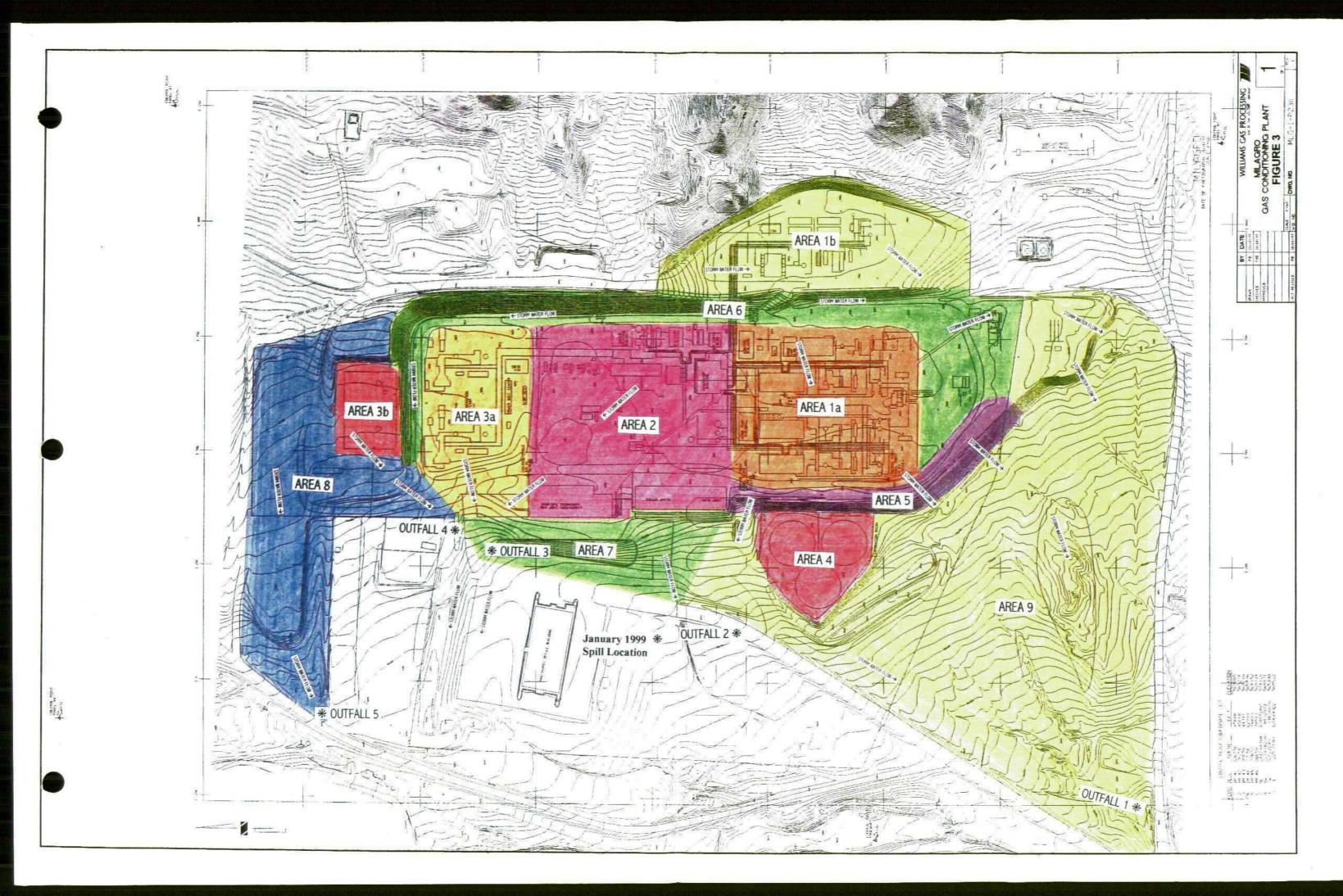
Figures





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# **SWP3 Documentation**

| If new, enter generic permit, otherwise enter previous permit:       Image: Substance in the provide permit.         B. Facility Operator Information       1. Name:       Substance in the provide permit.       2. Phone:       Substance in the provide permit.         3. Mailing Address:       a. Street or P.O. Box:       Image: Substance in the permit   | PDES<br>Form<br>510-6  | United States Environmental Protection Agency<br>Washington, DC 20460<br>Notice of Intent for Storm Water Discharges Assoc<br>INDUSTRIAL ACTIVITY Under the Multi-sector NPDES C   | OMB No. 2040-0086  |
|---|--|--|--|
| If new, enter generic permit, otherwise enter previous permit;       Image: Sector A         B. Facility Operator Information       1. Name: MITLELLIAMS IFIELD SERVIJ GES       2. Phone: SOSI/32/4/bit         3. Mailing Address: a. Street or PO. Box:       US EIS       2. Phone: SOS/6/32/4/bit         b. City:       BLoDEmFileLO       c. StateMIM       d. Zip Code:       SIJI413         1. Facility/Site Information       1. Facility/Site Information         1. Facility/Site Information       1. Facility/Site Information         1. Facility/Site Information       1. Facility/Site Information         1. Facility/Site Information       c. County: Sain Jule And         J. Coration Address:       a. Street:       U.L. Claid Claid         b. City:       BLIDOMFILE/       c. County: Sain Jule And         d. State:       File Address       a. Permit Applicant:       G. Congitude:         d. a. Permit Applicant:       Federal       State       Tuble G       Other public entity         b. Is the facility located on Indian Country Lands?       Yes       No       Stote A       Stote A         S. Does the facility discharge storm water rinc:       a. Receiving water(s)?       Wife Sain Call (SC) Codes or the 2-letter Advity Codes that best represent the principal products produced or services rendered by your facility and major co-located activities:         Primary:  | discharge polluta<br>ater Multi-sector<br>ection B of this for<br>thall applicable to<br>maintaining elig<br>ays after a compompleted. Please  | ants to waters of the United States, from the facility or site identified in Se<br>General Permit (MSGP). Submission of the NOI also constitutes notice<br>orm has read, understands, and meets the eligibility conditions of Part I of the<br>erms and conditions of the MSGP; understands that continued authorization<br>pibility for coverage, and that implementation of the permittee's pollution pre-<br>polete NOI is mailed. In order to be granted coverage, all information rece<br>e read and make sure you comply with all permit requirements, including the | ection C, under EPA's Storm<br>that the party identified in<br>he MSGP; agrees to comply<br>under the MSGP is contigent<br>evention plan is required two<br>guired on this form must be  |
| B. Facility Operator Information 1. Name: MITILLIIAMS IF IELD SERVIIGES 2. Phone: E05543244 3. Mailing Address: a. Street or P.O. Box: LISS FIR 14900 5. City: BL002mFileL0 1. C. Facility/Site Information 1. Facility/Site Information 2. Sector A Sector K Sector N Sector X Sector |  |  | W Permit Number(EPA Use Only)  |
| 1.Facility/Site Name:       IIIIIAA6R0       6.AS       IPROCESSI (IV)       ILIAAT         2.Location Address:       a. Street:       IIIIIIA       CIRINERSI (IV)       C. County:       San Julan         b. City:       BilliomAilel       IIIII       C. County:       San Julan       G. County:       San Julan         d. State:       MIM       e. Zip Code:       B'IIIII       F. Latitude:       g. Longitude:       G. County:         3.If you are filing as a co-permittee, enter storm water general permit number:       g. Longitude:       G. County:       San Julan         4.a. Permit Applicant:       [Federal]       [State]       Tribal       [Private]       Other public entity         b. Is the facility located on Indian Country Lands?       [Yes       [No       Score A       Score A         5.Does the facility located on Indian Country Lands?       [Yes       [No       IIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIII  | 1. Name: NIJI<br>3. Mailing Addre  | LLIIAMS FIELD SERVIJGES 2. Pho<br>ess: a. Street or P.O. Box: 188 ER 4900  |  |
| Do you certify under penalty of law that this document and all attachments were prepared under your direction supervision in accordance with a system designed to assure that qualified personnel properly gather and eval information submitted? Based on your inquiry of the person or persons who manage the system, or those pedirectly responsible for gathering the information, do you certify that the information submitted is, to the best knowledge and belief, true, accurate, and complete? Do you certify that you are aware that there are significated for submitting false information, including the possibility of fine and imprisonment for knowing violation.  | <ol> <li>Facility/Site N</li> <li>Location Addi<br/>b. City: B</li> <li>State: MI</li> <li>State: MI</li> <li>If you are filin</li> <li>A. Permit App</li> <li>b. Is the facili</li> <li>Does the facili</li> <li>Does the facili</li> <li>Does the facili</li> <li>Boes the facili</li> <li>A municipal<br/>If yes, nam</li> <li>The 4-digit S<br/>principal proop<br/>Primary: I</li> <li>Applicable set<br/>of the MSGP<br/>covered under<br/>Sector B</li> <li>Sector C</li> <li>Sector D</li> </ol> | Iame: <b>HILGAGROLOGAS PROGESSIPHOLOGAS</b> ress:       a. Street: <b>LAS CIR 14909</b> inses:       a. Street: <b>LAS CIR 14909 DomAiler CIR 14909 I I DomAiler G Sain Jun M</b> e. Zip Code: <b>B17141 B C</b> . County: <b>Sain M</b> e. Zip Code: <b>B17141 f</b> Latitude:       III <b>g</b> Longit <b>M</b> e. Zip Code: <b>B17141 f</b> Latitude:       IIII       g. Longit <b>M</b> e. Zip Code: <b>B17141 f</b> Latitude:       IIIIIII       g. Longit <b>M</b> e. Zip Code: <b>B17141 f</b> Latitude:       IIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIIII              | that best represent the activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activities:<br>activit |
|   | Do you certify<br>supervision ir<br>information s<br>directly respo<br>knowledge at<br>penalties for<br>Print Name: b  | y under penalty of law that this document and all attachments were prepare<br>in accordance with a system designed to assure that qualified personnel pro-<br>ubmitted? Based on your inquiry of the person or persons who manage to<br>possible for gathering the information, do you certify that the information su<br>and belief, true, accurate, and complete? Do you certify that you are aware<br>submitting false information, including the possibility of fine and imprison  | roperly gather and evaluate t<br>he system, or those persons<br>ibmitted is, to the best of you<br>e that there are significant<br>iment for knowing violations?   |
|   | Signature:   |  | Date: Dili2999<br>Page 1   |

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| To obtain a copy of the EPA's storm water multi-sector permit terms and conditions to which you are now held accountable, please call the EPA Office of Water Resource Center at (202) 260-7786. If you have general questions concerning the storm water program, |
|--|

| Ň   | NON-STORM WATER DISCHARGE   | CHARGE   | Worksheet #5<br>Completed by:  |  |  |
|---|---|--|--|--|--|
| ASSI  | ASSESSMENT AND CERTIFICATION  | FICATION   |  | Senior Environmental Specialist<br>7/22/2001   | and the second |
| Date of   |   | Method I lsed to   | be Results fro   |  | Name of Person Who   |
| Test or<br>Evaluation   | Outfall Directly Observed<br>During the Test (identify as<br>indicated on the site map)   | Test or Evaluate<br>Discharge  | Presence of Non-Storm Water<br>Discharge   | Identify Potential Significant<br>Sources  | Conducted the Test or<br>Evaluation  |
|   |   |  |  |  | Mark J. Bareta   |
|   |   |  |  |  |  |
|   |   |  |  |  |  |
|   |   |  |  |  |  |
|   |   |  |  |  |  |
|   |   |  |  |  |  |
|   |   |  | CERTIFICATION  |  |  |
| ا, <u>کولترمم</u><br>under my dire<br>submitted. Bi<br>information su<br>information ir | I, Jeffront Davance Corport<br>under my direction or supervision in accordance with a syster<br>submitted. Based on my inquiry of the person or persons who<br>information submitted is, to the best of my knowledge and bel<br>information. including the possibility of fine and imprisonment | (responsible corpore<br>dance with a system<br>rson or persons who<br>knowledge and belie<br>and imprisonment fo | I, Jeffrory Douvers (responsible corporate official), certify under penalty of law that this document and all attachments were prepared under my direction or supervision in accordance with a system designed to assure that qualified personnel properly gather and evaluate the information submitted. Based on my inquiry of the person or persons who manage the system or those persons directly responsible for gathering the information, the information submitted is, to the best of my knowledge and belief, true, accurate, and complete. I am aware that there are significant penalties for submitting false information, including the possibility of fine and imprisonment for knowing violations. | that this document and all atta<br>nnel properly gather and evalua<br>irectly responsible for gathering<br>ware that there are significant p | chments were prepared<br>ate the information<br>the information, the<br>benalties for submitting false           |
| A. Name & C   | A. Name & Official Title (type or print)  | Feer Phomere   | T) reference   | B. Area Code and Telephone No.<br>デット 632 ならっ」   | e No.  |
| C. Signature  |   | 2  |  | D. Date Signed   |  |
|   |   |  |  |  |  |

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Milagro Gas Processing Plant

# Appendix B

# **Inspection and Monitoring Worksheets**

#### Quarterly Inspection Report Milagro Gas Conditioning Plant

Inspection Date: \_\_\_\_\_

Name of Inspector: \_\_\_\_\_\_ Title: \_\_\_\_\_\_

Pollutant Reduction Measures Evaluation:

| Structural and<br>Sediment Controls | Physical Condition of<br>Control Features | Required Actions                       |
|-------------------------------------|---|--|
| Berms                               |   |  |
| Concrete Containment                |   |  |
| Gravel and Paving                   |   |  |
| Rip-Rap and Filter Fabric           |   |  |
| Ditches                             |   |  |
| Culverts and Piping                 |   |  |
| Ponds                               |   | ······································ |

#### **Best Management Practices Evaluation:**

| Best Management Practice                    | Condition of Practice | Required Actions |
|---|-----------------------|------------------|
| Good Housekeeping                           |                       |                  |
| Preventative Maintenance                    |                       |                  |
| Spill Prevention and Response<br>Procedures |                       |                  |
| Sediment and Erosion Control                |                       |                  |
| Process and Materials Control               |                       |                  |
| Process and Materials Management            |                       |                  |
| Employee Training                           |                       |                  |

#### **Additional Comments:**

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# Quarterly Storm Water Visual Monitoring Report Milagro Gas Conditioning Plant

Quarter: Year:

Inspection Date:

Name of Inspector:

Title:

Nature of Discharge: (Rain or Snowmelt)

| utfall 1<br>utfall 2<br>utfall 3<br>utfall 4 | Odor | Clarity /<br>Opacity | Floating<br>Solids | Clarity / Floating Suspended<br>Opacity Solids Solids | Settled<br>Solids | Foam | Oil Sheen | Foam Oil Sheen Comments |
|--|------|----------------------|--------------------|---|-------------------|------|-----------|-------------------------|
|  |      |                      |                    |   |                   |      |           |                         |
|  |      |                      |                    |   |                   |      |           |                         |
|  |      |                      |                    |   |                   |      |           |                         |
|  |      |                      |                    |   |                   |      |           |                         |
|  |      |                      |                    |   |                   |      |           |                         |

Comments:

#### Annual Comprehensive Compliance Evaluation Report Milagro Gas Conditioning Plant

Inspection Date: \_\_\_\_\_

Name of Inspector: \_\_\_\_\_ Title: \_\_\_\_\_

Others Present:

**Pollutant Source Evaluation:** 

| Pollutant Source Area                            | Pollutant Sources Observed | Required Actions |
|--|----------------------------|------------------|
| Plant Process Area Gas Tranes 1 - 4<br>(Area 1a) |                            |                  |
| Plant Process Area Gas Trane 5<br>(Area 1b)      |                            |                  |
| Utility Area<br>(Area 2)                         |                            | ····             |
| Cogeneration Area<br>(Area 3a)                   |                            |                  |
| Switch Yard<br>(Area 3b)                         |                            |                  |
| Wastewater Evaporation Pond Area (Area 4)        |                            |                  |
| Slope Drainage Area<br>(Area 5)                  |                            |                  |
| South Embankment Area<br>(Area 6)                |                            |                  |
| Plant Parking Area<br>(Area 7)                   |                            |                  |
| North 40 Lay-Down Area<br>(Area 8)               |                            |                  |
| South Discharge Area<br>(Area 9)                 |                            |                  |

**Comments:** 

Comprehensive Compliance Evaluation Report

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**Pollutant Reduction Measures Evaluation:** 

| Structural and<br>Sediment Controls | Physical Condition of<br>Control Features | Required Actions |
|-------------------------------------|---|------------------|
| Berms                               |   |                  |
| Concrete Containment                |   |                  |
| Gravel and Paving                   |   |                  |
| Rip-Rap and Filter Fabric           |   |                  |
| Ditches                             |   |                  |
| Culverts and Piping                 |   |                  |
| Detention Ponds                     |   |                  |
|                                     |   |                  |

#### **Best Management Practices Evaluation:**

| Best Management Practice                    | Condition of Practice | Required Actions                         |
|---|-----------------------|--|
| Good Housekeeping                           |                       | na an a |
| Preventative Maintenance                    |                       |  |
| Spill Prevention and Response<br>Procedures |                       |  |
| Sediment and Erosion Control                |                       |  |
| Process and Materials Control               |                       | <u></u>                                  |
| Process and Materials Management            |                       |  |
| Employee Training                           |                       |  |
|   |                       |  |

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#### **Additional Comments:**

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Notes: Supplementary documentation to the inspection report may include field notebooks, timed and dated photographs or videotapes, drawings, sketches, and maps.

Based on the comprehensive evaluation, the Multi-Sector General Permit (MSGP) requires that any potential pollutant sources or pollution prevention measures that are not adequately addressed in the Storm Water Pollution Prevention Plan (SWP3) need to be revised within 2 weeks of the evaluation. Any identified issues or measures that are revised in the SWMP must be implemented within 12 weeks after the evaluation as specified in the MSGP.

If no incidents of non-compliance are identified during the evaluation, the following certification of compliance with the SWP3 must be signed by the Plant Manager or other authorized individual within Williams Field Services organization. This authorization should be filed with the SWP3.

Certification of Compliance

I certify under penalty of law that this Comprehensive Compliance Evaluation Report was prepared under my direction or supervision and that the Milagro Gas Conditioning Plant is in compliance with the MSGP as specified in the SWP3. Based on my inquiry of the persons who conducted the evaluation, and those responsible for implementing the required actions, the information submitted is, to the best of my knowledge and belief, true, accurate, and complete.

Name:

Signature:

Date:

## Annual Comprehensive Compliance Evaluation Report Milagro Gas Conditioning Plant

Inspection Date: 6/26/2001

Name of Inspector: Clara Garcia Title: Environmental Compliance Administrator

Others Present: Eric Edmanson, Greg Millikin, J.D. Brow

**Pollutant Source Evaluation:** 

| Pollutant Source Area                            | Pollutant Sources Observed                   | Required Actions |
|--|--|------------------|
| Plant Process Area Gas Tranes 1 - 4<br>(Area 1a) | ОК   | None             |
| Plant Process Area Gas Trane 5<br>(Area 1b)      | ОК   | None             |
| Utility Area<br>(Area 2)                         | Steam Condensate and<br>D.I. Water discharge | . Stop Discharge |
| Cogeneration Area<br>(Area 3a)                   | ОК   | None             |
| Switch Yard<br>(Area 3b)                         | ОК   | None             |
| Wastewater Evaporation Pond Area<br>(Area 4)     | ОК   | None             |
| Slope Drainage Area<br>(Area 5)                  | ОК   | None             |
| South Embankment Area<br>(Area 6)                | ОК   | None             |
| Plant Parking Area<br>(Area 7)                   | See Area 2                                   | See Area 2       |
| North 40 Lay-Down Area<br>(Area 8)               | ок   | None             |
| South Discharge Area<br>(Area 9)                 | ОК   | None             |

#### **Comments:**

Steam condensate line was observed discharging outside the containment area. The condensate line will be redirected to discharge into the environmental drain system.

Evidence D.I. water discharge was observed outside the Water Treatment Building. D.I. water was discharged during filter renewal operations. D.I. water will be reused and not be discharged.

**Pollutant Reduction Measures Evaluation:** 

| Structural and<br>Sediment Controls | Physical Condition of<br>Control Features | Required Actions |
|-------------------------------------|---|------------------|
| Berms                               | ОК  | None             |
| Concrete Containment                | ОК  | None             |
| Gravel and Paving                   | ОК  | . None           |
| Rip-Rap and Filter Fabric           | ОК  | None             |
| Ditches                             | ОК  | None             |
| Culverts and Piping                 | ОК  | None             |
| Ponds                               | ОК  | None             |

Best Management Practices Evaluation:

| Best Management Practice                    | Condition of Practice | Required Actions |  |  |
|---|-----------------------|------------------|--|--|
| Good Housekeeping                           | ОК                    | None             |  |  |
| Preventative Maintenance                    | ОК                    | None             |  |  |
| Spill Prevention and Response<br>Procedures | ОК                    | None             |  |  |
| Sediment and Erosion Control                | ОК                    | None             |  |  |
| Process and Materials Control               | ОК                    | None             |  |  |
| Process and Materials Management            | ОК                    | None             |  |  |
| Employee Training                           | ОК                    | None             |  |  |

Additional Comments:

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Notes: Supplementary documentation to the inspection report may include field notebooks, timed and dated photographs or videotapes, drawings, sketches, and maps.

Based on the comprehensive evaluation, the Multi-Sector General Permit (MSGP) requires that any potential pollutant sources or pollution prevention measures that are not adequately addressed in the Storm Water Pollution Prevention Plan (SWP3) need to be revised within 2 weeks of the evaluation. Any identified issues or measures that are revised in the SWMP must be implemented within 12 weeks after the evaluation as specified in the MSGP.

If no incidents of non-compliance are identified during the evaluation, the following certification of compliance with the SWP3 must be signed by the Plant Manager or other authorized individual within Williams Field Services organization. This authorization should be filed with the SWP3.

#### Certification of Compliance

I certify under penalty of law that this Comprehensive Compliance Evaluation Report was prepared under my direction or supervision and that the Milagro Gas Conditioning Plant is in compliance with the MSGP as specified in the SWP3. Based on my inquiry of the persons who conducted the evaluation, and those responsible for implementing the required actions, the information submitted is, to the best of my knowledge and belief, true, accurate, and complete.

Name: Clara M Garcia Signature: Uaram. Bauel Date: August 13,2001

# WEEKLY CONTAINMENT INSPECTION REPORT MILAGRO GAS CONDITIONING PLANT

Inspection Date: 9-5-0/Name of Inspector: 10 Blow Title: Maint. Tech

Pollutant Reduction Measures Evaluation:

| Structural and<br>Sediment Controls | Physical Condition of<br>Control Features | <b>Required Actions</b> |
|-------------------------------------|---|-------------------------|
| Berms                               | oK  | Nonies                  |
| Concrete Containment                | OK  | NONE                    |
| Gravel and Paving                   | DK  | NOWE                    |
| Rip-Rap and Filter Fabric           | OK  | NONE                    |
| Ditches                             | ØK  | NONE                    |
| Culverts and Piping                 | 0K  | NANE                    |
| Ponds                               | DK  | NIDNE                   |
| Good Housekeeping                   | OK  | NONE                    |
| Preventative Maintenance            | NK  | NONE                    |

11

11

## Additional

Comments:

# Quarterly Inspection Report Milagro Gas Conditioning Plant

9-5-01 Inspection Date: D BROW \_\_\_\_\_ Title: MAINT. TECH. Name of Inspector:

**Pollutant Reduction Measures Evaluation:** 

| Structural and<br>Sediment Controls | Physical Condition of<br>Control Features | . Required Actions |
|-------------------------------------|---|--------------------|
| Berms                               | OK  | NONE               |
| Concrete Containment                | DK  | NANE               |
| Gravel and Paving                   | DK  | NONE               |
| Rip-Rap and Filter Fabric           | DK  | NONE               |
| Ditches                             | OK  | NONE               |
| Culverts and Piping                 | OK  | NONE               |
| Ponds                               | 0K  | NOVE               |

#### **Best Management Practices Evaluation:**

| Best Management Practice                    | Condition of Practice | Required Actions |
|---|-----------------------|------------------|
| Good Housekeeping                           | NK                    | None             |
| Preventative Maintenance                    | 0K                    | Nono             |
| Spill Prevention and Response<br>Procedures | DK.                   | NONE             |
| Sediment and Erosion Control                | DK.                   | NONE             |
| Process and Materials Control               | ÐK                    | NONE             |
| Process and Materials Management            | DK                    | NONE             |
| Employee Training                           | 0K                    | NONE             |

**Additional Comments:** 

Quarterly Storm Water Visual Monitoring Report Milagro Gas Conditioning Plant

Quarter: 340 Year: 200/

Inspection Date: 7-11-01

SPOL Ś Name of Inspector: \_\_\_\_

Title: Maint Tech

Nature of Discharge: (Rain or Snowmelt)

|  |   | 2            | <u>ې</u>     |                      | L            |
|--|---|--------------|--------------|----------------------|--------------|
| Foam Oil Sheen Comments  | X   | NO DISCHARGE | NO D'SCHARGE | X                    | No Discharge |
| Oil Sheen  | NONG  |              |              | NONE                 |              |
|  | NON   |              |              | NONE                 |              |
| Settled<br>Solids  | =na,N                                       |              |              | TAN MONE NONE OK     |              |
| Unusual Clarity / Floating Suspended<br>Odor Opacity Solids Solids | VEAR NOWE CLEAR NONE NONE NONE NONE NONE OK |              |              |                      |              |
| Floating<br>Solids   | None  |              |              | Nonë                 |              |
| Clarity /<br>Opacity   | CLORK                                       |              |              | NONE CLEDR NONE NONE |              |
| Unusual<br>Odor  | NONE  |              |              | NONE                 |              |
| Color  | Clerk                                       |              |              | ClEAR                |              |
| Date and<br>time   | 10-11-1                                     | 7-11.01      | 10-11-2      | 7-11-01              | 10-11-2      |
| Outfall Observation<br>Location                                    | Outfall 1                                   | Outfall 2    | Outfall 3    | Outfall 4            | Outfall 5    |

**Comments:** 

# Appendix C

# Storm Water Visual Monitoring Procedure

## Storm Water Visual Monitoring Procedure

#### **Outfalls and Drainage Points**

The Storm Water Pollution Prevention Plan (SWP3) includes a section that describes the outfall locations and the area drained by each outfall. A site drainage map provides a visual reference for the location of the outfalls and the outline of the drainage area to each outfall. Visual inspection and monitoring (if required) must usually be performed at all outfalls. If characteristics of storm water discharge are expected to be similar for different outfalls, then a facility may choose to report the test results from one outfall as representative of all outfalls.

#### Sheet Flow Run-off

Sheet flow run-off occurs in areas that have no discernible discrete discharge point. Vacant land and open parking lots are examples of sheet flow run-off areas. These areas should be described in the SWP3 and shown on the site drainage map.

#### Regulatory Requirements for Visual Inspection and Monitoring of Storm Water Discharges

Each facility covered under the multi-sector general storm water permit must conduct quarterly visual monitoring of storm water discharges at each outfall during qualifying storm events and must document the results.

In accordance with the general permit, visual examinations of storm water quality must be performed and documented at least once each calendar quarter. The visual inspections of storm water discharge at each outfall must be completed during a qualifying storm event during daylight hours. A qualifying storm event is defined as rainfall of at least 0.1 inches, with no previous storm events having occurred within the previous 72 hours (3-days). Realistically, however, based on conditions present at typical industrial sites, a rainfall rate of 0.3 to 0.5 inches is often required before an appreciable amount of flow occurs that can be collected for a representative observation. A rain gauge mounted in an accessible area outside, away from traffic and the possibility of disturbance, should be used to record the total amount of rainfall at the site over the duration of the storm event (e.g. 1.2 inches over 4.5 hours).

Visual inspections should be made within 30 minutes to 1 hour after the beginning of storm water flow through the outfalls. The inspections shall document observations for the following discharge characteristics:

- Color
- Unusual odor
- Clarity/opacity
- Floating, suspended, or settled solids
- Foam
- Oil sheen
- Other obvious indications of contamination or pollution (e.g. residues or foreign matter)

The examinations must be conducted in a well-lit area. No analytical testing is required for the quarterly visual monitoring requirements beyond the recording of physical observations. Documented results must be retained in the SWP3.

#### Resources/Coordination Needed to Conduct Visual Storm Water Monitoring

Storm water quarterly visual monitoring can be documented on a worksheet-type checklist with fields for entering comments. The worksheet can be formatted to include physical inspections of plant areas drained by each outfall, so that potential sources of contamination can be identified if such evidence is observed in the discharge. Take into consideration appropriate safety precautions and other logistics regarding outfall access during inspection planning. The following items are recommended to conduct the visual inspection at each outfall, or to gather additional data:

- Mounted rain gauge located outside in an accessible location of the plant,
- Watch or clock,

- Umbrella or rain-protective clothing,
- Skid resistant rubber or plastic safety boots,
- Clean, unused white plastic bucket with handle and plastic lid, at least 2-gallon capacity,
- Chemical resistant rope,
- 1-quart or equivalent unused clear glass sample jar or clear plastic container,
- Blank quarterly monitoring form.

The bucket(s) and jar(s) used for visual inspection of storm water discharges should be clean and reserved for this purpose only. The goal is to capture a representative volume of storm water at each outfall and record the observations made regarding physical appearance and characteristics of the discharge. The run off should be observed during the first thirty minutes to one hour after flow at the outfall begins. Once a storm event occurs, the period before a sufficient amount of flow is discharged can be as little as ten minutes or as much as an hour, depending on factors such as how far from the drainage area the actual outfall discharge point is.

Steps To Complete The Visual Inspection

1. Once the storm event begins, note the time. Check the flow at the closest outfall approximately every 15 minutes. The flow must be sufficient to collect at least half a bucket within a minute or two (approximately 1 gal/min). If sufficient flow is never achieved after the first hour, then the storm event will be insufficient for observation.

2. If sufficient flow is achieved within the first thirty minutes to one hour, record the date and beginning time that the outfall observations are made in the appropriate field on the worksheet form. Proceed to the first outfall.

3. Note the physical appearance of any run-on from adjacent property, if such is present. If the appearance of storm water coming from adjacent property appears to be different from that being discharged from the site, note such differences on the visual observation worksheet.

4. Except where noticeable unusual odors of a chemical nature are detected, it is presumed that storm water discharges may have a characteristic "earthy" or "moldy" smell, especially if there is decayed vegetation or algae growth in the discharge area. Some discharges may have no odor at all. It is helpful to record a baseline observation of the typical characteristics of an "unpolluted" discharge for reference. Unusual odors should be recorded on the worksheet, otherwise it is presumed that no unusual odor was present.

5. Collect at least half a bucket of run-off for visual inspection. Tie the rope to the bucket handle if needed to maneuver the bucket into the drainage point. Catch the flow in the most rapid portion of the stream without scraping dirt or debris from the ground.

6. As soon as a sufficient volume of run off is collected in the bucket, fill the sample jar with some of the collected run off. Keep the remaining water in the bucket, cover the container, and set aside temporarily.

7. Note the following parameters and quality of the discharge in the sample jar and record on the worksheet:

Color: Should normally be able to see through the water. Presence of dirt or excessive silt may indicate soil erosion or washout, or presence of other types of residues.

Excess floating, settled, and suspended solids: Solids may be present in the collected water, and these solids may float on the water, settle out in the bottom of the container, or remain suspended in the water. The source of these solids may be dirt, material washed out from a stockpile, construction area, etc., depending on the color. Normally these should not prevent seeing through the liquid in the jar. Note the presence of suspended and settleable solids by color, opacity of the liquid, and nature of the material if known.

8. Note the following parameters and quality of discharge from inspecting the water in the sample jar and comparing with the run-off left in the bucket, recording observations on the worksheet:

Foam: Presence of foam suggests that detergent residues, emulsifiers, or surfactants may have been deposited on pavements from illicit washing of vehicles or equipment outside. Also look at the outfall for signs of foaming.

Oil sheen: Use a flashlight if needed to observe the surface of the collected run-off in the bucket. Visible oil sheen indicates that petroleum or other oil-like residues have washed off pavement. In many cases, this can be attributed to oil staining from vehicles in parking lots, docks, or traffic areas, or from outside storage of containers that may be leaking or which have residues on the outside.

Excess settled solids and debris: Use flashlight if needed to look at collected run-off in the bucket and sample jar. Have solids settled at the bottom in a layer more than 0.25 inches? Are there any unidentified pieces of debris or trash other than grass, pebbles, or dirt?

9. Note the following parameters by observing the condition of the outfall drainage point and recording on the worksheet:

Staining: Does the outfall bear evidence of unusual staining other than that attributable to rust from metal storm grates or other identified sources?

Evidence of chemical dumping: Is there any evidence of debris associated with chemical residues or container remnants? Is there a tar-like discolored layer or film, or does the drainage area appear pitted or worn away from exposure to a corrosive material?

Dry Weather Flow: Prior to the storm event, was there any noticeable dry weather flow? If so, was the source determined? If yes, record this data on the worksheet.

10. After observations are complete at the first outfall, proceed to the next outfall or sheet flow run-off area to make similar observations, following Steps 1 - 9. If the storm event ceases prior to completing observations at all outfalls and sheet flow areas, note this on the worksheet. A separate qualifying storm event during the same quarter will be needed to complete observations for remaining outfalls or sheet flow areas.

After the visual inspection is completed, rinse the bucket and sample jar with potable water and stored dry in a secure area until next storm event.

11. After the storm event has ceased, record the total rainfall amount received at the mounted rain gauge at the facility, and the duration of the storm event (as best determined) on the worksheet. After collecting the rainfall data, empty the rain gauge for the next storm event.

Notes on collection of rainfall data: The rain gauge must be checked and emptied after each storm event, whether visual inspections are conducted or not, to ensure the accuracy of the data collected for the next storm event. An on-site rain gauge is generally more accurate than reports from nearby weather stations, as rainfall amounts may vary widely even over short distances.

The same worksheet used to record visual storm water quality inspection results may also be used to record physical plant area inspections, either before or after the qualifying storm event. After the form is completed, turn in to the SWP3 Team Coordinator for review; follow up action if needed, and archiving in the SWP3.

The following conditions should be recorded on a quarterly monitoring form if they occur:

- Lack of a qualifying storm event during a calendar quarter,
- Inability to conduct visual observations due to facility shut down,
- Occurrence of storm event when facility is unoccupied,
- Inability to conduct visual observations because storm event begins after daylight hours,
- Inability to conduct visual observation because of lack of sufficient flow, and
- Storm event occurs within three days of previous rainfall.

# SPE LTTR. DATED February 25, 1997 February 25, 1997 The Subsurface Investigation of The Milagro Plant Septic Leach Field San Juan Basin, New Mexico

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January 1997

**Prepared For** 

# WILLIAMS FIELD SERVICES SALT LAKE CITY, UTAH

Project 17309



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# TABLE OF CONTENTS

| 1.0 INTRODUCTION   | 1      |
|--|--------|
| 2.0 SCOPE OF WORK  | 1      |
| <ul><li>2.1 Subsurface Investigation</li><li>2.2 Sample Septic Tank</li><li>2.3 Profile Laboratory Waste</li></ul> | 2      |
| 3.0 METHODOLOGY  | 2      |
| <ul><li>3.1 Subsurface Investigation</li><li>3.2 Sample Septic Tank</li></ul>                                      | 2<br>3 |
| 4.0 RESULTS  | 4      |
| <ul><li>4.1 Soil Samples</li><li>4.2 Sludge/Waste Water</li><li>4.3 Chemical Waste Profile</li></ul>               | 4      |
| 5.0 SUMMARY  | 5      |

# LIST OF TABLES

TABLE 1 - Soil Analytical ResultsTABLE 2 - Sludge Analytical ResultsTABLE 3 - Waste Water Analytical Results

# LIST OF FIGURES

FIGURE 1 - Boring Locations, Septic Lines and Septic Tank Locations FIGURE 2 - Milagro Plant As-Built Diagram

# LIST OF APPENDICES

APPENDIX A - Field Documentation APPENDIX B - Laboratory Reports APPENDIX C - Chemical Waste Profile

**APPENDIX D** - Data Evaluation



## **1.0 INTRODUCTION**

On December 12, 1996, Philip Environmental Services Corporation (Philip) initiated field work for a subsurface investigation of Williams Field Services (WFS) Milagro Plant Septic Leach Field. The Milagro Plant (the Plant) is located northeast of Bloomfield, New Mexico, in Township - 29 North, Range - 11 West, Section 11. This investigation was required based on the plant laboratories past practice of disposing chemical wastes into the septic system. Chemical wastes are no longer disposed of in this manner; they are currently stored on-site prior to disposal at a licensed facility.

Prior to commencing with the field work for this project, Philip developed a work plan detailing the method of investigation, and the analytical parameters for samples. The work plan was approved, with modifications, on September 9, 1996, by the New Mexico Oil Conservation Division (NMOCD). Philip completed the following scope of work for this project.

### 2.0 SCOPE OF WORK

As part of the Scope of Work for this project, Philip completed three separate tasks, detailed below, to determine the extent, if any, of contamination from the discharged chemical wastes within the septic leach field.

#### 2.1 Subsurface Investigation

One soil boring was drilled within the approximate center of the septic leach field. One soil sample was collected from within the septic leach field area and one soil sample was collected from approximately 10 feet beneath the bottom of the leach field. Both samples were submitted for laboratory analysis.

In addition, two soil borings were drilled on the estimated downgradient edge of the leach field. Soil samples were collected from each boring at approximately 10 feet beneath the bottom of the leach field. Both samples were submitted for laboratory analysis. Boring locations, septic lines, and the septic tank locations are presented in Figure 1.



#### 2.2 Sample Septic Tank

One sludge sample and one waste water sample was collected from the Plant septic tank, and both samples were submitted for laboratory analysis. The sludge sample was submitted for Hazardous Waste Identification analysis. The waste water sample was submitted for the New Mexico Water Quality Control Commission (NMWQCC) parameters of concern, as per the NMOCD's request.

### 2.3 Profile Laboratory Waste

In addition to the field investigation, the laboratory chemical wastes were profiled for disposal, and accepted by Laidlaw Environmental Services' Clive, Utah, facility for disposal.

## **3.0 METHODOLOGY**

All soil samples were submitted for total Resource Conservation and Recovery Act (RCRA) metals (Method SW 6010/7470), semi-volatile organics (Method SW 8270), and volatile organics (Method SW 8260). Septic tank sludge samples were submitted for Hazardous Waste Identification, including Toxicity Characteristic Leaching Procedure (TCLP), ignitability, corrosivity, and reactivity. Septic tank waste water samples were submitted for NMWQCC parameters of concern, which included metals (SW846 6010), volatile organics (SW846 8260), and semi-volatile organics (SW846 8270). All sample containers were labeled with the appropriate analysis, date and time of collection, sample number, sample location, and sample collector. All sample identification numbers and requested analysis were documented on a Chain-of-Custody form. Samples were placed on ice, and shipped via overnight delivery to Zenon Environmental Laboratories following strict Chain-of-Custody procedures. All field documentation forms are presented in Appendix A.

#### 3.1 Subsurface Investigation

One soil boring (BH-1) was advanced in the approximate center of the leach field using a CME-75 drill rig and 4 1/4-inch inside diameter hollow-stem augers. Soil samples were collected at 5 foot intervals, using a split-spoon soil sampler. The samples were screened for volatiles with a photoionization detector (PID). No volatiles were detected with the



PID. One soil sample was collected from approximately 3 - 5 feet below ground surface (bgs) for laboratory analysis. This sample appeared to exhibit the most moisture, indicating the sample was collected from within the discharge area of the leach field. A second sample was collected from approximately 10 - 12 feet bgs and submitted for laboratory analysis. The boring was then grouted to the surface with a neat cement slurry containing 5 percent bentonite. A description of soils encountered in each boring was recorded on individual Record of Subsurface Exploration forms, found in Appendix A.

A second boring (BH-2) was advanced on the approximate downgradient side of the leach field. Soil samples were collected at 5 foot intervals, and were screened for volatiles using a PID. No volatiles were detected with the PID. One soil sample was collected from approximately 15 - 17 feet bgs for laboratory analysis. The boring was then grouted to the surface in the same manner as BH-1.

A third boring (BH-3) was advanced to 15 feet bgs. BH-3 is located southwest of BH-1, on the outside edge of the leach field. Soil samples were collected at 5 foot intervals, and were screened for volatiles using a PID. No volatiles were detected with the PID. Refusal was encountered in cobbles at approximately 13 feet bgs, and the soil sample from 13 - 15 feet bgs was submitted for laboratory analysis. The boring was then grouted to the surface in the same manner as BH-1.

One background sample was collected on September 23, 1996, using a stainless-steel hand auger. The sample was collected topographically upgradient from the septic tank. The hand auger was advanced to approximately 5 feet bgs, and a soil sample was collected for laboratory analysis. A separate aliquot of soil was collected at this depth and screened for volatiles. No volatiles were detected with the PID. The boring was then backfilled with native soil.

#### 3.2 Sample Septic Tank

On January 24, 1997, Philip sampled the septic tank associated with the Plant laboratory. Two samples were collected from the tank (sludge, and waste water) and submitted for laboratory analysis. The sludge was collected using a steel grab sampler which was decontaminated with a potable water and Alconox<sup>TM</sup> soap wash, followed by a distilled



water rinse, prior to use. The waste water was collected using a disposable polyethylene bailer.

## 4.0 **RESULTS**

The leach field soil analytical results are presented in Table 1, with the laboratory report presented in Appendix B. The septic tank sludge and waste water analytical results are presented in Table 2 and Table 3, respectively, with the laboratory report presented in Appendix B.

#### 4.1 Soil Samples

With the exception of barium, all metals were at or below method detection limits (MDL). In all samples, including the background sample, acetone, and methylene chloride, were detected at levels slightly above the MDL. Toluene was detected in BH-1 (3-5) at the MDL. After evaluation of the laboratory report, the presence of acetone, methylene chloride, and toluene, were attributed to laboratory interference. Documentation of the data evaluation is presented in Appendix D. All semi-volatiles were at or below the MDL for each parameter. In addition, groundwater was not encountered at any of the boring locations.

#### 4.2 Sludge/Waste Water

The septic tank sludge was analyzed for Hazardous Waste Identification, including TCLP, ignitability, corrosivity, and reactivity. The TCLP volatiles, semi-volatiles, and metals analyses were all below regulatory limits. In addition, results for ignitability, corrosivity, and reactivity were within regulatory limits, which classifies the septic sludge as non-hazardous.

The septic waste water was analyzed for NMWQCC parameters of concern. All of the parameters with the exception of phenol were below NMWQCC levels. Phenol was reported at 11 micrograms per Liter ( $\mu$ g/L), slightly above the NMWQCC limit of 5  $\mu$ g/L. Based on the analytical results for the leach field soil samples, it does not appear that the soil was impacted by phenols. Several non-NMWQCC parameters were detected at levels above the corresponding MDL. Benzoic acid, which is not listed under NMWQCC and



does not appear to be listed as a Hazardous Waste, was detected at  $620 \mu g/L$ . Based on telephone discussions with the RCRA hotline, benzoic acid by itself does not meet the requirements for toxicity. The Plant laboratory personnel indicated they were not aware of using any chemicals containing benzoic acid. Based on the analytical results for the leach field soil samples, it does not appear that the soil has been impacted by benzoic acid.

## 4.3 Chemical Waste Profile

Used laboratory chemical wastes were profiled, and accepted for disposal at a permitted Laidlaw Environmental Services' facility. The waste was profiled as RCRA Non Hazardous/Exempt. The Waste Profile sheet is presented in Appendix C. The chemical wastes will be lab-packed prior to transporting, and will be picked up from the laboratory for disposal at Laidlaw's facility on an as needed basis.

## 5.0 SUMMARY

The laboratory analytical results indicate minimal impact to the soils from discharged used chemical wastes, within and downgradient of the septic leach field. Analytical results for the septic tank sludge/waste water also indicate minimal impact from the discharged chemical wastes. The local geology indicates shallow bedrock (sandstone), as documented in the boring log for BH-2. Also, groundwater was not encountered in any boring within or downgradient of the septic leach field. On average, depth to groundwater in the vicinity of the Plant ranges from 30 feet to 100 feet bgs ("Hydrogeology and Water Resources of San Juan Basin, New Mexico" 1983, & "Availability of Hydrologic Data in San Juan County, New Mexico", 1984). In addition, laboratory chemical wastes are no longer disposed of into the septic leach field, and are currently stored on-site prior to disposal at a permitted facility. Based on these factors, Philip believes no further action is warranted at this site.



**TABLE 1 - SOIL ANALYTICAL RESULTS**

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MILAGRO PLANT SEPTIC FIELD

#### SOIL ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

|           | Ē                 | PART I - M          | ETALS                  |                     |                        |      |
|-----------|-------------------|---------------------|------------------------|---------------------|------------------------|------|
| PARAMETER | <b>BH-1</b> (3-5) | <b>BH-1</b> (10-12) | <b>BH-2</b><br>(15-17) | <b>BH-3</b> (13-15) | <b>BG-1</b> *<br>(5-7) | MDL  |
|           | Te                | OTAL RCRA           | METALS                 |                     |                        |      |
| Arsenic   | <10               | <10                 | <10                    | <10                 | 2.2                    | 5.0  |
| Barium    | 280               | 130                 | 190                    | 150                 | 140                    | 0.1  |
| Cadmium   | <0.4              | <0.4                | <0.4                   | <0.4                | <0.4                   | 0.2  |
| Chromium  | <10               | <10                 | <10                    | <10                 | <10                    | 5    |
| Lead      | <20               | <20                 | <20                    | <20                 | <20                    | 10   |
| Mercury   | <0.04             | < 0.04              | < 0.04                 | < 0.04              | <0.04                  | 0.04 |
| Selenium  | <20               | <20                 | <20                    | <20                 | <0.5                   | 10   |
| Silver    | <1.0              | <1.0                | <1.0                   | <1.0                | <1.0                   | 0.5  |

All results reported in milligrams per kilogram (mg/kg)

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

< - Less than MDL

The second se

and the second s

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\* - BG-1 collected on 9/23/96

#### (Continued)

## SOIL ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

| and the second | BH-1   | BH-1       | BH-2    | BH-3    | BG-1  |       |
|--|--------|------------|---------|---------|-------|-------|
| PARAMETER  | (3-5)  | (10-12)    | (15-17) | (13-15) | (5-7) | MDL   |
|  | ······ | OLATILE OI |         | L       | l     |       |
| Acetone  | 0.18   | 0.15       | 0.14    | 0.12    | 0.072 | 0.030 |
| Acrolein   | <      | <          | <       | <       | <     | 0.010 |
| Acrylonitrile  | <      | <          | <       | <       | <     | 0.010 |
| Benzene  | <      | <          | <       | <       | <     | 0.005 |
| Bromoform  | <      | <          | <       | <       | <     | 0.010 |
| Bromomethane   | <      | <          | <       | <       | <     | 0.010 |
| 2-Butanone   | <      | <          | <       | <       | <     | 0.015 |
| Carbon Disulfide   | <      | <          | <       | <       | <     | 0.010 |
| Carbon Tetrachloride   | <      | <          | <       | <       | <     | 0.010 |
| Chlorobenzene  | <      | <          | <       | <       | <     | 0.005 |
| Chlorodibromomethane   | <      | <          | <       | <       | <     | 0.005 |
| Chloroethane   | <      | <          | <       | <       | <     | 0.010 |
| 2-Chloroethylvinylether  | <      | <          | <       | <       | <     | 0.010 |
| Chloroform   | <      | <          | <       | <       | <     | 0.005 |
| Chloromethane  | <      | <          | <       | <       | <     | 0.010 |
| 1,2-Dichlorobenzene  | <      | <          | <       | <       | <     | 0.005 |
| 1,3-Dichlorobenzene  | <      | <          | <       | <       | <     | 0.005 |
| 1,4-Dichlorobenzene  | 0.005  | <          | <       | <       | 0.008 | 0.005 |
| Dichlorobromomethane   | <      | <          | <       | <       | <     | 0.005 |
| 1,1-Dichloroethane   | <      | <          | <       | <       | <     | 0.005 |
| 1,2-Dichloroethane   | <      | <          | <       | <       | <     | 0.005 |
| 1,1-Dichloroethene   | <      | <          | <       | <       | <     | 0.010 |
| cis-1,2-Dichloroethene   | <      | <          | <       | <       | <     | 0.010 |
| trans-1,2-Dichloroethene   | <      | <          | <       | <       | <     | 0.010 |
| 1,2-Dichloropropane  | <      | <          | <       | <       | <     | 0.005 |
| cis-1,3-Dichloropropene  | <      | <          | <       | <       | <     | 0.005 |
| trans-1,3-Dichloropropene  | <      | <          | <       | <       | <     | 0.005 |
| Ethylbenzene   | <      | <          | <       | <       | <     | 0.005 |
| 2-Hexanone   | <      | <          | <       | <       | <     | 0.010 |

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

< - Less than MDL

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#### (Continued)

## SOIL ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

| PART II - VOLATILE ORGANICS<br>(Continued)   |                   |                        |                        |                     |                      |           |  |
|--|-------------------|------------------------|------------------------|---------------------|----------------------|-----------|--|
| PARAMETER  | <b>BH-1</b> (3-5) | <b>BH-1</b><br>(10-12) | <b>BH-2</b><br>(15-17) | <b>BH-3</b> (13-15) | <b>BG-1</b><br>(5-7) | MDL       |  |
|  | N N               | OLATILE OI             | RGANICS                |                     | <u> </u>             | 1 <u></u> |  |
| Methylene Chloride   | 0.039             | 0.034                  | 0.29                   | 0.033               | 0.091                | 0.010     |  |
| 4-Methyl-2-Pentanone   | <                 | <                      | <                      | <                   | <                    | 0.010     |  |
| Styrene  | <                 | <                      | <                      | <                   | <                    | 0.005     |  |
| 1,1,1,2-Tetrachloroethane  | <                 | <                      | <                      | <                   | <                    | 0.010     |  |
| 1,1,2,2-Tetrachloroethane  | <                 | <                      | <                      | <                   | <                    | 0.010     |  |
| Tetrachloroethene  | <                 | <                      | <                      | <                   | <                    | 0.020     |  |
| Toluene  | 0.005             | <                      | <                      | <                   | 0.005                | 0.005     |  |
| 1,1,1-Trichloroethane  | <                 | <                      | <                      | <                   | <                    | 0.005     |  |
| 1,1,2-Trichloroethane  | <                 | <                      | <                      | <                   | <                    | 0.010     |  |
| Trichloroethene  | <                 | <                      | <                      | <                   | <                    | 0.005     |  |
| Vinyl Acetate  | <                 | <                      | <                      | <                   | <                    | 0.010     |  |
| Vinyl Chloride   | <                 | <                      | <                      | <                   | <                    | 0.010     |  |
| Xylenes (Total)  | <                 | <                      | <                      | <                   | 0.005                | 0.005     |  |
| All results reported in milligram<br>MDL - Method Detection Limit<br>< - Less than MDL |                   |                        | er Sheet for M         | DL comments         | )                    |           |  |

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#### SOIL ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

|                             | PART III - S |             |         |         |        |      |
|-----------------------------|--------------|-------------|---------|---------|--------|------|
|                             | BH-1         | BH-1        | BH-2    | BH-3    | BG-1   |      |
| PARAMETER                   | (5-7)        | (25-27)     | (17-18) | (20-22) | (5-7)  | MDL  |
| ·                           |              | II-VOLATILE |         |         |        |      |
| Phenol                      | <0.22        | < 0.22      | < 0.22  | < 0.22  | < 0.22 | 0.11 |
| Bis(2-chloroethyl)ether     | < 0.36       | < 0.36      | < 0.36  | < 0.36  | < 0.36 | 0.18 |
| 2-Chlorophenol              | <0.54        | <0.54       | <0.54   | <0.54   | <0.54  | 0.27 |
| 1,3-Dichlorobenzene         | <0.40        | <0.40       | <0.40   | < 0.40  | < 0.40 | 0.20 |
| 1,4-Dichlorobenzene         | <0.40        | < 0.40      | < 0.40  | <0.40   | < 0.40 | 0.20 |
| 1,2-Dichlorobenzene         | <0.40        | < 0.40      | < 0.40  | <0.40   | <0.40  | 0.20 |
| Bis(2-chloroisopropyl)ether | < 0.30       | < 0.30      | < 0.30  | < 0.30  | < 0.30 | 0.15 |
| Hexachloroethane            | <0.40        | <0.40       | <0.40   | <0.40   | <0.40  | 0.20 |
| N-Nitrosodi-N-Propylamine   | <0.42        | <0.42       | < 0.42  | < 0.42  | < 0.42 | 0.21 |
| Nitrobenzene                | <0.40        | < 0.40      | < 0.40  | < 0.40  | <0.40  | 0.20 |
| Isophorone                  | < 0.80       | <0.80       | < 0.80  | < 0.80  | < 0.80 | 0.40 |
| 2-Nitrophenol               | < 0.28       | < 0.28      | <0.28   | <0.28   | < 0.28 | 0.14 |
| 2,4-Dimethylphenol          | < 0.34       | < 0.34      | < 0.34  | < 0.34  | < 0.34 | 0.17 |
| Bis(2-chloroethoxy)methane  | < 0.26       | < 0.26      | < 0.26  | < 0.26  | < 0.26 | 0.13 |
| 2,4-Dichlorophenol          | < 0.24       | < 0.24      | < 0.24  | < 0.24  | < 0.24 | 0.12 |
| 1,2,4-Trichlorobenzene      | <0.40        | <0.40       | <0.40   | < 0.40  | < 0.40 | 0.20 |
| Naphthalene                 | < 0.06       | <0.06       | <0.06   | < 0.06  | < 0.06 | 0.03 |
| Hexachlorobutadiene         | <0.40        | <0.40       | <0.40   | < 0.40  | < 0.40 | 0.20 |
| 4-Chloro-3-Methylphenol     | <0.28        | < 0.28      | < 0.28  | < 0.28  | <0.28  | 0.14 |
| Hexachlorocyclopentadiene   | < 0.40       | <0.40       | <0.40   | <0.40   | < 0.40 | 0.20 |
| 2,4,6-Trichlorophenol       | <0.24        | < 0.24      | <0.24   | < 0.24  | < 0.24 | 0.12 |
| 2-Chloronaphthalene         | <0.18        | <0.18       | < 0.18  | < 0.18  | < 0.18 | 0.09 |
| Acenaphthylene              | < 0.08       | < 0.08      | < 0.08  | < 0.08  | < 0.08 | 0.04 |
| Dimethyl phthalate          | < 0.22       | < 0.22      | < 0.22  | < 0.22  | < 0.22 | 0.11 |
| 2,6-Dinitrotoluene          | <0.12        | <0.12       | < 0.12  | <0.12   | < 0.12 | 0.06 |
| Acenaphthene                | <0.14        | < 0.14      | <0.14   | <0.14   | < 0.14 | 0.07 |
| 2,4-Dinitrophenol           | <0.96        | <0.96       | < 0.96  | <0.96   | < 0.96 | 0.48 |
| 2,4-Dinitrotoluene          | <0.10        | <0.10       | <0.10   | <0.10   | < 0.10 | 0.05 |
| 4-Nitrophenol               | < 0.28       | < 0.28      | <0.28   | <0.28   | < 0.28 | 0.14 |
| Fluorene                    | < 0.06       | < 0.06      | < 0.06  | <0.06   | < 0.06 | 0.03 |
| 4-Chlorophenylphenylether   | < 0.18       | <0.18       | <0.18   | <0.18   | <0.18  | 0.09 |
| Diethyl phthalate           | < 0.22       | < 0.22      | < 0.22  | < 0.22  | < 0.22 | 0.11 |
| 4,6-Dinitro-2-methylphenol  | < 0.30       | < 0.30      | < 0.30  | < 0.30  | < 0.30 | 0.15 |
| N-Nitrosodiphenylamine      | < 0.38       | < 0.38      | < 0.38  | < 0.38  | < 0.38 | 0.19 |

All results reported in milligrams per kilogram (mg/kg)

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

< - Less than MDL

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### SOIL ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

|                            | BH-1   | BH-1        | BH-2     | BH-3    | BG-1   |      |
|----------------------------|--------|-------------|----------|---------|--------|------|
| PARAMETER                  | (5-7)  | (25-27)     | (17-18)  | (20-22) | (5-7)  | MDL  |
| · · · · ·                  | SEN    | 41-VOLATILE | ORGANICS |         |        |      |
| 4-Bromophenylphenylether   | < 0.06 | <0.06       | <0.06    | <0.06   | <0.06  | 0.03 |
| Hexachlorobenzene          | <0.40  | <0.40       | <0.40    | <0.40   | <0.40  | 0.20 |
| Pentachlorophenol          | <0.22  | <0.22       | <0.22    | <0.22   | <0.22  | 0.11 |
| Phenanthrene               | < 0.06 | <0.06       | <0.06    | < 0.06  | <0.06  | 0.03 |
| Anathracene                | <0.04  | <0.04       | < 0.04   | < 0.04  | < 0.04 | 0.02 |
| Di-n-butyl phthalate       | < 0.22 | < 0.22      | < 0.22   | < 0.22  | <0.22  | 0.11 |
| Fluoranthene               | <0.04  | <0.04       | <0.04    | < 0.04  | < 0.04 | 0.02 |
| Pyrene                     | <0.06  | <0.06       | < 0.06   | < 0.06  | <0.06  | 0.03 |
| Benzyl butyl phthalate     | <0.12  | <0.12       | <0.12    | <0.12   | < 0.12 | 0.06 |
| 3,3-Dichlorobenzidine      | < 0.20 | < 0.20      | < 0.20   | < 0.20  | < 0.20 | 0.10 |
| Benzo(a)anthracene         | <0.04  | <0.04       | < 0.04   | < 0.04  | < 0.04 | 0.02 |
| Chrysene                   | <0.06  | <0.06       | <0.06    | < 0.06  | <0.06  | 0.03 |
| Bis(2-ethylhexyl)phthalate | <0.28  | <0.28       | <0.28    | < 0.28  | <0.28  | 0.14 |
| Di-n-octyl phthalate       | <0.22  | <0.22       | <0.22    | < 0.22  | < 0.22 | 0.11 |
| Benzo(b)fluoranthene       | <0.08  | <0.08       | <0.08    | < 0.08  | < 0.08 | 0.04 |
| Benzo(k)fluoranthene       | <0.08  | <0.08       | <0.08    | <0.08   | < 0.08 | 0.04 |
| Benzo(a)pyrene             | <0.10  | <0.10       | <0.10    | <0.10   | <0.10  | 0.05 |
| Indeno(1,2,3-cd)pyrene     | <0.12  | <0.12       | <0.12    | <0.12   | <0.12  | 0.06 |
| Dibenzo(a,h)anthracene     | < 0.08 | < 0.08      | < 0.08   | < 0.08  | < 0.08 | 0.04 |
| Benzo(g,h,i)perylene       | <0.08  | < 0.08      | <0.08    | < 0.08  | <0.08  | 0.04 |
| N-Nitrosodimethylamine     | <2.0   | <2.0        | <2.0     | <2.0    | <2.0   | 1.0  |
| Aniline                    | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| Benzyl alcohol             | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| Carbazole                  | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| 2-Methylphenol             | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| Benzoic acid               | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| 4-Chloroaniline            | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| 2-Methylnaphthalene        | < 0.20 | <0.20       | <0.20    | <0.20   | <0.20  | 0.10 |
| 2,4,5-Trichlorophenol      | < 0.20 | < 0.20      | <0.20    | <0.20   | <0.20  | 0.10 |
| 2-Nitroaniline             | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| 3-Nitroaniline             | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| Dibenzofuran               | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| Benzidine                  | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |
| 4-Nitroaniline             | <1.0   | <1.0        | <1.0     | <1.0    | <1.0   | 0.50 |

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**TABLE 2 - SLUDGE ANALYTICAL RESULTS**

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MILAGRO PLANT SEPTIC TANK

#### SLUDGE ANALYTICAL RESULTS MILAGRO SEPTIC TANK

| PART I - TCLP METALS |                       |                |       |
|----------------------|-----------------------|----------------|-------|
| PARAMETER            | STS-12497<br>(Sludge) | TCLP<br>LIMITS | MDL   |
| Mercury              | <                     | 0.2            | 0.05  |
| Arsenic              | <0.55                 | 5.0            | 0.50  |
| Barium               | 1.2                   | 100.0          | 0.001 |
| Cadmium              | 0.002                 | 1.0            | 0.002 |
| Chromium             | <                     | 5.0            | 0.004 |
| Lead                 | <0.022                | 5.0            | 0.020 |
| Selenium             | <0.066                | 1.0            | 0.060 |
| Silver               | < 0.011               | 5.0            | 0.010 |

All results reported in milligrams per liter (mg/L)

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments) < - Less than MDL

TCLP Metals analyzed using SW 846 Method 6010 except Mercury

Mercury analyzed via SW 846 Method 7470

#### (Continued)

## SLUDGE ANALYTICAL RESULTS MILAGRO SEPTIC TANK

|                      | STS-12497 | TCLP   |     |
|----------------------|-----------|--------|-----|
| PARAMETER            | (Sludge)  | LIMITS | MDL |
| Benzene              | <2.0      | 500    | 0.2 |
| 2-Butanone           | <50       | NA     | 5.0 |
| Carbon Tetrachloride | <3.0      | 500    | 0.3 |
| Chlorobenzene        | <6.0      | 10000  | 0.6 |
| Chloroform           | <4.0      | 6000   | 0.4 |
| 1,2-Dichloroethane   | <4.0      | 500    | 0.4 |
| 1,1-Dichloroethene   | <4.0      | 700    | 0.4 |
| Pyridine             | <2500     | 5000   | 250 |
| Tetrachloroethene    | <2.0      | NA     | 0.2 |
| Trichloroethene      | <3.0      | NA     | 0.3 |
| Vinyl Chloride       | <29       | 200    | 2.9 |
| 1,4-Dichlorobenzene  | 13        | 7500   | 0.2 |

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

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#### SLUDGE ANALYTICAL RESULTS MILAGRO SEPTIC TANK

| PART III - TCLP SEMI-VOLATILE ORGANICS |                       |                |     |  |
|--|-----------------------|----------------|-----|--|
| PARAMETER                              | STS-12497<br>(Sludge) | TCLP<br>LIMITS | MDL |  |
| o-Cresol                               | <                     | 200000         | 1.7 |  |
| m & p-Cresol                           | <                     | 200000         | 3.5 |  |
| 1,4-Dichlorobenzene                    | 7.4                   | 7500           | 2.0 |  |
| 2,4-Dinitrotoluene                     | <                     | 130            | 0.5 |  |
| Nitrobenzene                           | <                     | 2000           | 2.0 |  |
| Pentachlorophenol                      | <                     | 100000         | 1.1 |  |
| 2,4,5-Trichlorophenol                  | <                     | 400000         | 0.6 |  |
| 2,4,6-Trichlorophenol                  | <                     | 2000           | 1.2 |  |
| Hexachloroethane                       | <                     | 3000           | 2.0 |  |
| Hexachlorobutadiene                    | <                     | 500            | 2.0 |  |
| Hexachlorobenzene                      | <                     | 130            | 2.0 |  |

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

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### SLUDGE ANALYTICAL RESULTS MILAGRO SEPTIC TANK

| PARAMETER                                 | Units  | STS-12497<br>(Sludge) |
|---|--------|-----------------------|
| Corrosivity                               | pН     | 7.85                  |
| Reactivity (reaction/flume/dust/gas)      | Deg C  | 0                     |
| Reactivity (temperature change 10 min)    | Deg C  | 0                     |
| Reactivity (temperature change - initial) | Deg C  | 0                     |
| Ignitability                              | mm/min | <                     |
| pH analyzed via SW 846 9045               |        |                       |
| Reactivity analyzed via ASTMD 5058C       |        |                       |

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 TABLE 3 - WASTE WATER ANALYTICAL RESULTS

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MILAGRO PLANT SEPTIC TANK

#### WASTE WATER ANALYTICAL RESULTS MILAGRO SEPTIC TANK

| PART I - METALS |                     |                  |        |
|-----------------|---------------------|------------------|--------|
| PARAMETER       | ST-12497<br>(Water) | NMWQCC<br>LIMITS | MDL    |
| Arsenic         | <                   | 0.1              | 0.0020 |
| Selenium        | <                   | 0.05             | 0.0020 |
| Aluminum        | 0.26                | NA               | 0.030  |
| Barium          | 0.49                | 1.0              | 0.001  |
| Beryllium       | <                   | NA               | 0.001  |
| Cadmium         | 0.004               | 0.01             | 0.002  |
| Calcium         | 39                  | NA               | 0.20   |
| Chromium        | <                   | 0.05             | 0.004  |
| Cobalt          | < 0.011             | 0.05             | 0.010  |
| Copper          | 0.11                | 1.0              | 0.005  |
| Iron            | 0.38                | 1.0              | 0.010  |
| Lead            | < 0.022             | 0.05             | 0.020  |
| Magnesium       | 7.5                 | NA               | 0.050  |
| Manganese       | 0.031               | 0.2              | 0.006  |
| Molybdenum      | 0.017               | 1.0              | 0.010  |
| Nickel          | < 0.011             | 0.2              | 0.010  |
| Phosphorus      | 8.6                 | NA               | 0.060  |
| Potassium       | 23                  | NA               | 1.00   |
| Silver          | < 0.011             | 0.05             | 0.010  |
| Sodium          | 44                  | NA               | 0.10   |
| Thallium        | <0.066              | NA               | 0.060  |
| Vanadium        | < 0.006             | NA               | 0.005  |
| Zinc            | 0.21                | 10.0             | 0.005  |

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments) < - Less than MDL

Metals analyzed using SW 846 Method 6010 except Arsenic and Selenium

Arsenic analyzed via SW 846 Method 7060

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Selenium analyzed via SW 846 Method 7740

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## WASTE WATER ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

| PART II - VOLATILE ORGANICS |                     |                        |       |
|-----------------------------|---------------------|------------------------|-------|
| PARAMETER                   | ST-12497<br>(Water) | NMWQCC<br>LIMITS(µG/L) | MDL   |
| Acetone                     | 14                  | NA                     | 0.030 |
| Acrolein                    | <                   | NA                     | 0.010 |
| Acrylonitrile               | <                   | NA                     | 0.010 |
| Benzene                     | <                   | 10.0                   | 0.005 |
| Bromoform                   | <                   | NA                     | 0.010 |
| Bromomethane                | <                   | NA                     | 0.010 |
| 2-Butanone                  | <                   | NA                     | 0.015 |
| Carbon Disulfide            | <                   | NA                     | 0.010 |
| Carbon Tetrachloride        | <                   | 10.0                   | 0.010 |
| Chlorobenzene               | <                   | NA                     | 0.005 |
| Chlorodibromomethane        | <                   | NA                     | 0.005 |
| Chloroethane                | <                   | NA                     | 0.010 |
| 2-Chloroethylvinylether     | <                   | NA                     | 0.010 |
| Chloroform                  | 4.8                 | 100.0                  | 0.005 |
| Chloromethane               | <                   | NA                     | 0.010 |
| 1,2-Dichlorobenzene         | <                   | NA                     | 0.005 |
| 1,3-Dichlorobenzene         | <                   | NA                     | 0.005 |
| l,4-Dichlorobenzene         | 37                  | NA                     | 0.005 |
| Dichlorobromomethane        | <                   | NA                     | 0.005 |
| 1,1-Dichloroethane          | <                   | 25.0                   | 0.005 |
| 1,2-Dichloroethane          | <                   | 10.0                   | 0.005 |
| 1,1-Dichloroethene          | <                   | NA                     | 0.010 |
| cis-1,2-Dichloroethene      | <                   | NA                     | 0.010 |
| rans-1,2-Dichloroethene     | <                   | NA                     | 0.010 |
| 1,2-Dichloropropane         | <                   | NA                     | 0.005 |
| cis-1,3-Dichloropropene     | <                   | NA                     | 0.005 |
| trans-1,3-Dichloropropene   | <                   | NA                     | 0.005 |
| Ethylbenzene                | <                   | 750.0                  | 0.005 |
| 2-Hexanone                  | <                   | NA                     | 0.010 |

All results reported in micrograms per liter ( $\mu$ g/L)

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MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

< - Less than MDL

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#### WASTE WATER ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

| PART II - VOLATILE ORGANICS<br>(Continued) |                     |                        |     |  |
|--|---------------------|------------------------|-----|--|
| PARAMETER                                  | ST-12497<br>(Water) | NMWQCC<br>LIMITS(µg/L) | MDL |  |
| Methylene Chloride                         | 39                  | 100.0                  | 2.0 |  |
| 4-Methyl-2-Pentanone                       | <                   | NA                     | 2.0 |  |
| Styrene                                    | <                   | NA                     | 1.0 |  |
| 1,1,1,2-Tetrachloroethane                  | <                   | NA                     | 2.0 |  |
| 1,1,2,2-Tetrachloroethane                  | <                   | 10.0                   | 2.0 |  |
| Tetrachloroethene                          | <                   | NA                     | 4.0 |  |
| Toluene                                    | <                   | 750.0                  | 1.0 |  |
| 1,1,1-Trichloroethane                      | <                   | 60.0                   | 1.0 |  |
| 1,1,2-Trichloroethane                      | <                   | 10.0                   | 2.0 |  |
| Trichloroethene                            | <                   | NA                     | 1.0 |  |
| Vinyl Acetate                              | <                   | NA                     | 2.0 |  |
| Vinyl Chloride                             | <                   | 1.0                    | 2.0 |  |
| Xylenes (Total)                            | <                   | 620.0                  | 1.0 |  |

All results reported in micrograms per liter ( $\mu$ g/L)

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

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#### WASTE WATER ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

| PART III - SEMI-VOLATILE ORGANICS ST-12497 NMWQCC |         |        |     |  |
|---|---------|--------|-----|--|
| PARAMETER   | (Water) | LIMITS | MDL |  |
| Phenol  | 11      | 5.0    | 1.1 |  |
| Bis(2-chloroethyl)ether                           | <       | NA     | 1.8 |  |
| 2-Chlorophenol                                    | <       | NA     | 2.7 |  |
| 1,3-Dichlorobenzene                               | <       | NA     | 2.0 |  |
| 1,4-Dichlorobenzene                               | 20      | NA     | 2.0 |  |
| 1,2-Dichlorobenzene                               | <       | NA     | 2.0 |  |
| Bis(2-chloroisopropyl)ether                       | <       | NA     | 1.5 |  |
| Hexachloroethane                                  | <       | NA     | 2.0 |  |
| N-Nitrosodi-N-Propylamine                         | <       | NA     | 2.1 |  |
| Nitrobenzene                                      | <       | NA     | 2.0 |  |
| Isophorone  | <       | NA     | 4.0 |  |
| 2-Nitrophenol                                     | <       | NA     | 1.4 |  |
| 2,4-Dimethylphenol                                |         | NA     | 1.7 |  |
| Bis(2-chloroethoxy)methane                        | <       | NA     | 1.3 |  |
| 2,4-Dichlorophenol                                | <       | NA     | 1.2 |  |
| 1,2,4-Trichlorobenzene                            | <       | NA     | 2.0 |  |
| Naphthalene                                       | <       | 30     | 0.3 |  |
| Hexachlorobutadiene                               | <       | NA     | 2.0 |  |
| 4-Chloro-3-Methylphenol                           | <       | NA     | 1.4 |  |
| Hexachlorocyclopentadiene                         | <       | NA     | 2.0 |  |
| 2,4,6-Trichlorophenol                             | <       | NA     | 1.2 |  |
| 2-Chloronaphthalene                               | <       | NA     | 0.9 |  |
| Acenaphthylene                                    | <       | NA     | 0.4 |  |
| Dimethyl phthalate                                | <       | NA     | 1.1 |  |
| 2,6-Dinitrotoluene                                | <       | NA     | 0.6 |  |
| Acenaphthene                                      | <       | NA     | 0.7 |  |
| 2,4-Dinitrophenol                                 | <       | NA     | 4.8 |  |
| 2,4-Dinitrotoluene                                | <       | NA     | 0.5 |  |
| 4-Nitrophenol                                     | <       | NA     | 1.4 |  |
| Fluorene  | <       | NA     | 0.3 |  |
| 4-Chlorophenylphenylether                         | <       | NA     | 0.9 |  |
| Diethyl phthalate                                 | <       | NA     | 1.1 |  |
| 4,6-Dinitro-2-methylphenol                        | <       | NA     | 1.5 |  |
| N-Nitrosodiphenylamine                            | <       | NA     | 1.9 |  |

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

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### WASTE WATER ANALYTICAL RESULTS MILAGRO SEPTIC FIELD

| PARAMETER                  | ST-12497<br>(Water) | NMWQCC<br>LIMITS | MDL |
|----------------------------|---------------------|------------------|-----|
| 4-Bromophenylphenylether   | <                   | NA               | 0.3 |
| Hexachlorobenzene          | <                   | NA               | 2.0 |
| Pentachlorophenol          | <                   | NA               | 1.1 |
| Phenanthrene               | <                   | NA               | 0.3 |
| Anathracene                | <                   | NA               | 0.2 |
| Di-n-butyl phthalate       | <                   | NA               | 1.1 |
| Fluoranthene               | <                   | NA               | 0.2 |
| Pyrene                     | <                   | NA               | 0.3 |
| Benzyl butyl phthalate     | 1.0                 | NA               | 0.6 |
| 3,3-Dichlorobenzidine      | <                   | NA               | 1.0 |
| Benzo(a)anthracene         | <                   | NA               | 0.2 |
| Chrysene                   | <                   | NA               | 0.3 |
| Bis(2-ethylhexyl)phthalate | 2.1                 | NA               | 1.4 |
| Di-n-octyl phthalate       | <                   | NA               | 1.1 |
| Benzo(b)fluoranthene       | <                   | NA               | 0.4 |
| Benzo(k)fluoranthene       | <                   | NA               | 0.4 |
| Benzo(a)pyrene             | <                   | 0.7              | 0.5 |
| Indeno(1,2,3-cd)pyrene     | <                   | NA               | 0.6 |
| Dibenzo(a,h)anthracene     | <                   | NA               | 0.4 |
| Benzo(g,h,i)perylene       | <                   | NA               | 0.4 |
| N-Nitrosodimethylamine     | <                   | NA               | 10  |
| Aniline                    | <                   | NA               | 5.0 |
| Benzyl alcohol             | <                   | NA               | 2.0 |
| Carbazole                  | <                   | NA               | 5.0 |
| 2-Methylphenol             | <                   | NA               | 1.7 |
| Benzoic acid               | 620                 | NA               | 1.0 |
| 4-Chloroaniline            | <                   | NA               | 5.0 |
| 2-Methylnaphthalene        | <                   | NA               | 1.0 |
| 2,4,5-Trichlorophenol      | <                   | NA               | 0.6 |
| 2-Nitroaniline             | <                   | NA               | 5.0 |
| 3-Nitroaniline             | <                   | NA               | 5.0 |
| Dibenzofuran               | <                   | NA               | 5.0 |
| Benzidine                  | <                   | NA               | 10  |
| 4-Nitroaniline             | <                   | NA               | 5.0 |

All results reported in micrograms per liter ( $\mu g/L$ ))

MDL - Method Detection Limit (See Laboratory Report Cover Sheet for MDL comments)

< - Less than MDL

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FIGURE 1 - BORING LOCATIONS, SEPTIC LINES AND SEPTIC TANK LOCATIONS

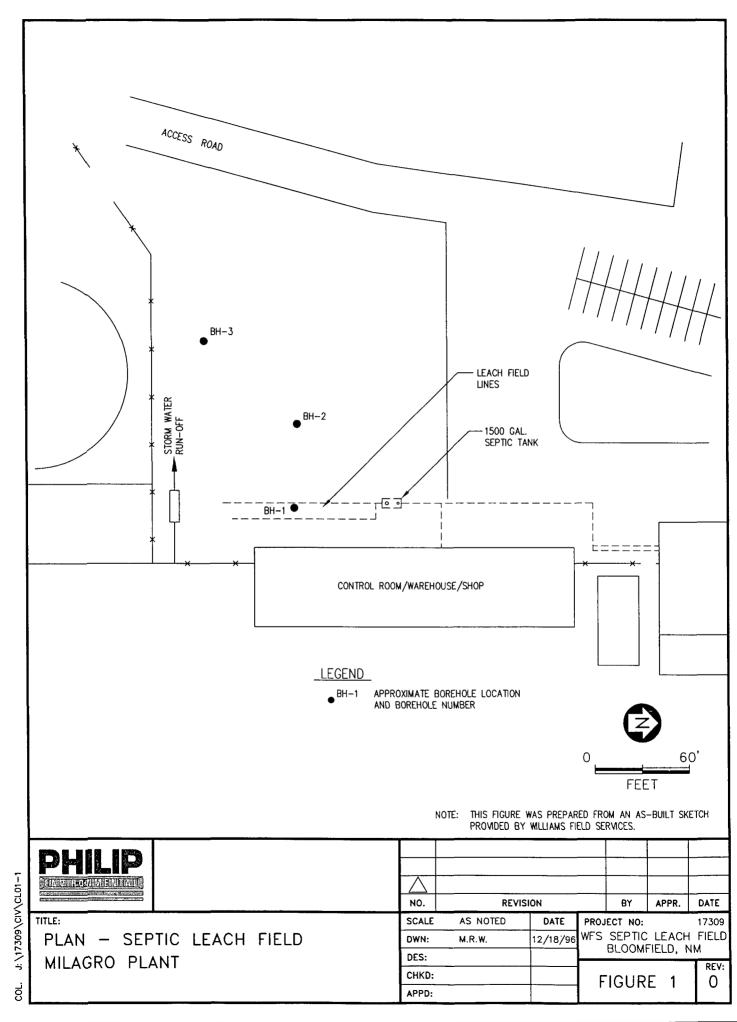
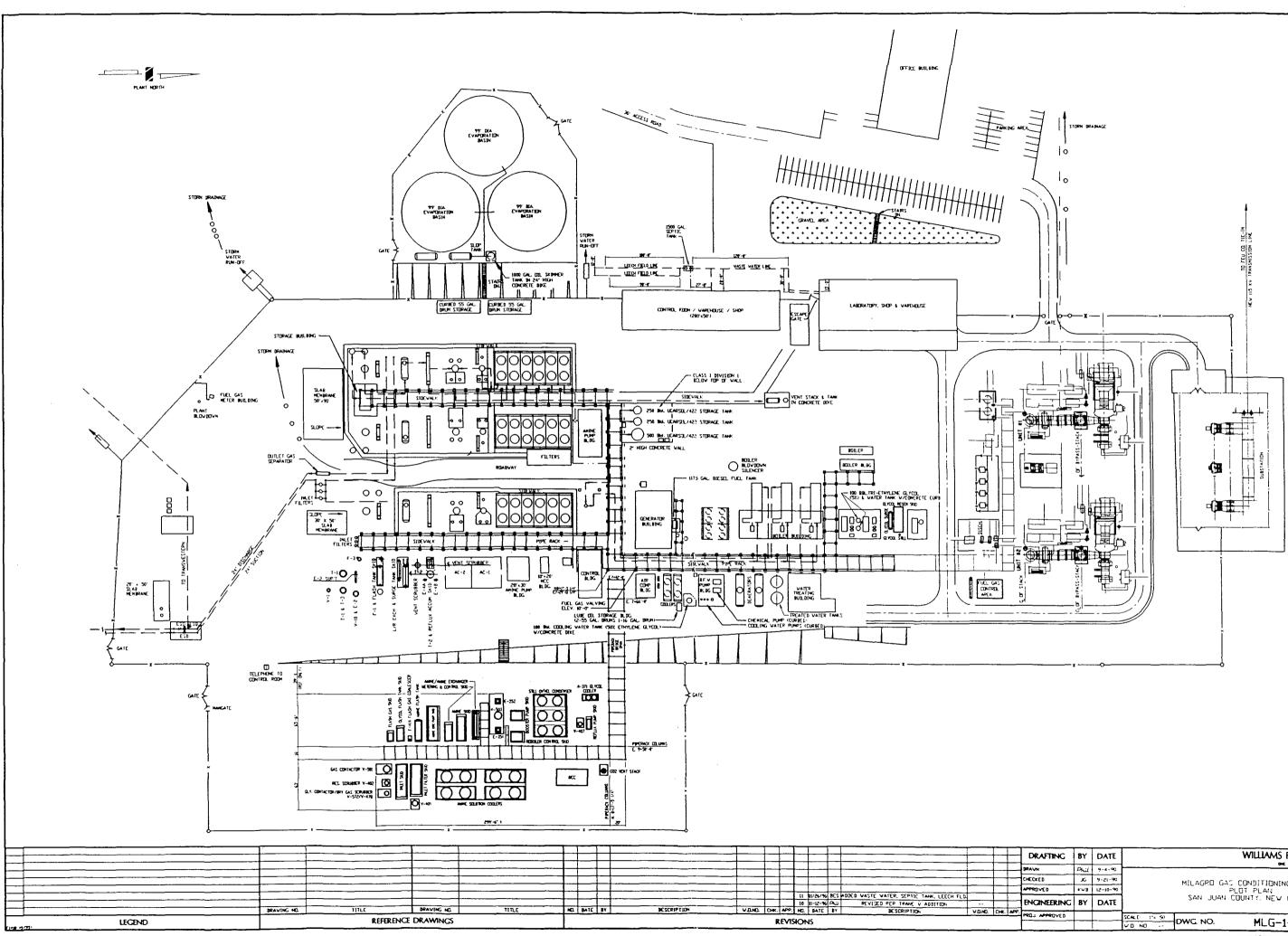


FIGURE 2 - MILAGRO PLANT AS-BUILT DIAGRAM

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|       |      |          |   | ALY      |
|-------|------|----------|---|----------|
| NG    | 8Y   | DATE     | WILLIAMS FIELD SERVICES                   |          |
|       | Phil | 9-4-90   |   | =        |
|       | ж    | 9-21-90  | MILAGRE GAS CONDITIONING PLANT            |          |
| _     | KVB  | 12-10-90 | PLOT PLAN                                 |          |
| RING  | BY   | DATE     | SAN JUAN COUNTY, NEW MEXICO               | 6CV#     |
| IVE D |      |          | SCALE 12= 50<br>V.D. ND DWG. NO. MLG-1-P1 | REV.     |
|       |      | <u> </u> |   | <u> </u> |

**APPENDIX A - FIELD DOCUMENTATION** 

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### **RECORD OF SUBSURFACE EXPLORATION**

PHILIP ENVIRONMENTAL SERVICES INC.

4000 Monroe Road Farmington, New Mexico 87401

(505) 326-2262 FAX (505) 326-2388

| Elevation         |                       |
|-------------------|-----------------------|
| Borehole Location | Center of Leach Field |
| GWL Depth         | N/A                   |
| Logged By         | CM Chance             |
| Drilled By        | K Padilla             |
| Date/Time Started | 12/12/96 - 0830       |
| Date/Time Comple  | ted 12/12/96 - 093    |

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|             | Borehole #     | BH- 1       |
|-------------|----------------|-------------|
|             | Welt #         |             |
|             | Page <u>1</u>  | of <u>1</u> |
|             |                |             |
| WFS Milagr  | 0              |             |
| 17039       | Phase          | 6001.77     |
| Milagro Pla | nt, Bloomfield | , NM        |
|             |                |             |
| CM (        | Chance         |             |
|             |                |             |

Personnel On-Site Contractors On-Site Client Personnel On-Site

Project Name

Project Number Project Location

Well Logged By

F. Rivera

nep

L. Gooding

Drilling Method 4 1/4 ID Hollow Stem Auger Air Monitoring Method PID

| Depth  | Sample | Sample   | Sample<br>Type & | Sample Description   | USCS   | Depth<br>Lithology |    | r Monitor |   | Drilling Conditions |
|--------|--------|----------|------------------|--|--------|--------------------|----|-----------|---|---------------------|
| (Feet) | Number | Interval | Recovery         | Classification System: USCS  | Symbol | Change             |    | Inits: PP |   | & Blow Counts       |
| 0      |        |          | (inches)         |  |        | (feet)             | BZ | ВН        | s |                     |
| 5      | 1      | 3-5      | 6                | Brown sandy Clay, medium plastic, soft<br>very moist   |        |                    | 0  | 0         | 0 | -0830 hour - Samp   |
| 10     | 2      | 10-12    | 24               | Brown siltly SAND, fine-medium sand,<br>loose, grades to clayey sand, trace<br>gypsum parting, dry |        |                    | 0  | 0         | 0 | -0839 hour - Samp   |
| 15<br> |        |          |                  | TOB 12'  |        |                    |    |           |   |                     |
| 20     |        |          |                  |  |        |                    |    |           |   |                     |
| 25     |        |          |                  |  |        |                    |    |           |   |                     |
| 30<br> |        |          |                  |  |        |                    |    |           |   |                     |
| 35     |        |          |                  |  |        |                    |    |           |   |                     |
| 40     |        |          |                  |  |        |                    |    |           |   |                     |

from 3-5' (BH1-3-5) and 10-12' (BH1-10-12) and submitted for laboratory analysis. Groundwater not encountered.

Geologist Signature

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### RECORD OF SUBSURFACE EXPLORATION

PHILIP ENVIRONMENTAL SERVICES INC.

4000 Monroe Road Farmington, New Mexico 87401 (505) 326-2262 FAX (505) 326-2388

| Elevation         |                      |
|-------------------|----------------------|
| Borehole Location | West of Leach Field  |
| GWL Depth         | N/A                  |
| Logged By         | CM Chance            |
| Drilled By        | K Padilla            |
| Date/Time Started | 12/12/96 - 0915      |
| Date/Time Comple  | eted 12/12/96 - 1000 |

|             | Borehole #     | BH- 2   |
|-------------|----------------|---------|
|             | Well #         |         |
|             | Page 1         | of 1    |
|             |                |         |
| WFS Milag   | ro             |         |
| 17039       | Phase          | 6001.77 |
| Milagro Pla | nt, Bloomfield | , NM    |
|             |                |         |
| CM          | Chance         |         |

Personnel On-Site Contractors On-Site Client Personnel On-Site

Well Logged By

Project Name

Project Number **Project Location** 

F. Rivera

L. Gooding

Drilling Method Air Monitoring Method

4 1/4 ID Hollow Stem Auger PID

| Depth<br>(Feet) | Sample<br>Number | Sample<br>Interval | Sample<br>Type &<br>Recovery<br>(inches) | Sample Description Classification System: USCS | USCS<br>Symbol | Depth<br>Lithology<br>Change<br>(feet) | 1 | r Monitor<br>Inits: PPI<br>BH |   | Drilling Conditions<br>& Blow Counts |
|-----------------|------------------|--------------------|--|--|----------------|--|---|-------------------------------|---|--------------------------------------|
|                 |                  |                    | Type &<br>Recovery<br>(inches)           |  |                | Lithology<br>Change                    | ι | Inits: PP                     | м |                                      |
| 30              |                  |                    |  |  |                |  |   |                               |   |                                      |

Comments:

Boring terminated at 17 feet BGS. Sample collected from 15-17 feet BGS (BH2-15-17) and submitted for laboratory analysis. Borehole grouted to the surface. Groundwater not encountered.

**Geologist Signature** 

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#### **RECORD OF SUBSURFACE EXPLORATION**

PHILIP ENVIRONMENTAL SERVICES INC.

4000 Monroe Road Farmington, New Mexico 87401 (505) 326-2262 FAX (505) 326-2388

| Elevation         |                     |
|-------------------|---------------------|
| Borehole Location | SW of Leach Field   |
| GWL Depth         | N/A                 |
| Logged By         | CM Chance           |
| Drilled By        | K Padilla           |
| Date/Time Started | 12/12/96 - 1015     |
| Date/Time Comple  | ted 12/12/96 - 1100 |

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|               | Borehole #    | BH- 3       |
|---------------|---------------|-------------|
|               | Well #        |             |
|               | Page <u>1</u> | of <u>1</u> |
| WFS Milagro   |               |             |
| 17039         | Phase         | 6001.77     |
| Milagro Plant | , Bloomfield, | NM          |

Well Logged By Personnel On-Site Contractors On-Site Client Personnel On-Site

Project Name

Project Number

Project Location

CM Chance F. Rivera

L. Gooding

Drilling Method 4

4 1/4 ID Hollow Stem Auger od PID

|                 |        |          | Sample               |  | l        | Depth            |        |                  |          |   |
|-----------------|--------|----------|----------------------|--|----------|------------------|--------|------------------|----------|---|
| Depth<br>(Feat) | Sample | Sample   | Type &               | Sample Description   | USCS     | Lithology        | 1      | r Monitor        |          | Drilling Conditions                         |
| (Feet)          | Number | Interval | Recovery<br>(inches) | Classification System: USCS  | Symbol   | Change<br>(feet) | BZ U   | Jnits: PPI<br>BH | M<br>S   | & Blow Counts                               |
| 0               |        |          |                      | Brown sandy CLAY, soft, medium<br>plastic, slightly moist                        |          |                  | 0      | 0                | 0        | -Cuttings                                   |
| 10              |        |          |                      | Light groutoff white sites CAND which it   |          |                  | ~      |                  |          | Cobbles @ 6'                                |
|                 | 1      | 13-15    |                      | Light grey/off white silty SAND, slightly<br>cemented, fine-v. fine sand, gravel |          |                  | 0      | 0                | 0        | Cobbles @ 10'<br>Refusal w/ augers<br>@ 13' |
| 15<br>20<br>25  |        |          |                      | As Above<br>TOB 15'  |          |                  | 0      | 0                | Ο        | -1046 hour - Sample                         |
|                 |        |          |                      |  |          |                  |        |                  |          |   |
| Comments:       |        | Boring   | terminat             | ed at 15 foot BGS. Sample collected from 13                                      | -15 feet | (BH3-13          | -151 8 |                  | ad for l | aboratory                                   |

analysis. Borehole grouted to surface. Groundwater not encountered.

iony Chance

|          |            | F   | hase   |   |  |
|----------|------------|---|--|---|--|
| Services |            |   | 11450  | .Task No. <u>60</u>   | 00.77  |
| VI. T. V |            |   |  |   |  |
| tic lann |            |   |  |   |  |
|          |            |   |  |   |  |
| Portabl  | e Screenin | ig Instr  | umen   | t Used  | Nor  |
|          | Туре       |   |  | Manufacture   | n Mode   |
|          | PID (Larr  | ιp  | eV)  |   |  |
| on 🛛     | FID        |   |  |   |  |
|          | CGI        |   |  |   |  |
| space    |            |   |  |   |  |
| ing U    | Other      |   |  |   |  |
|          | Other      |   |  |   |  |
|          |            |   |  |   | Field  |
|          |            |   |  |   | Instrumen  |
| Collect  | ed Sar     | nple ly   | /pe  | Collected   | Reading  |
|          | Soil       | Sed.  | Slg.   | l   |  |
| 1015     |            |   | $\checkmark$   | 1-250-P   | NA   |
| 1015     |            |   | $\checkmark$   | 1-250mp   | NA   |
|          | ĺ          |   |  |   |  |
|          |            |   |  |   |  |
|          |            |   |  |   |  |
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|          |            |   |  |   |  |
|          | 1          | 1 1   |  |   |  |
|          |            |   |  |   |  |
|          | Portabl    | Portable Screenin<br>Type<br>PID (Lan<br>PID (Lan<br>PID<br>CGI<br>Space<br>ing<br>Dther<br>Time<br>Collected<br>Soil<br>1015 | Portable Screening Instructive<br>Type<br>PID (Lamp<br>PID (Lamp<br>CG1<br>space<br>ing<br>Dther<br>Time<br>Collected Sample Ty<br>Soit Sed.<br>1015<br>1015 | Portable Screening Instrument<br>Type<br>PID (LampeV)<br>on<br>FID<br>CGI<br>space<br>ing<br>Other<br>Other<br>Time<br>Collected<br>Sample Type<br>Soit Sed. Sig.<br>IDIS<br>IDIS | Portable Screening Instrument Used         Type       Manufacture         Image: PID       < |

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| Sampler Water Quality Beadings Water Collection Data (Explain in Co  | PHI                   | LIF             | Locat         | tion No          |                    |               |              |             |                 |               | W                   | ΤΑ                    | ER S                                   | SAN         | /IPLII    |                                       |
|--|-----------------------|-----------------|---------------|------------------|--------------------|---------------|--------------|-------------|-----------------|---------------|---------------------|-----------------------|--|-------------|-----------|---------------------------------------|
| Sampling Specifications       Initial Measurements         Requested Sampling<br>Depth Interval (feet)   | ENVIRONM              | IENITA          |               |                  |                    | ,             |              |             |                 |               |                     |                       |  |             |           |                                       |
| Sampling Specifications       Initial Measurements         Requested Sampling  | ample Tvp             | be: [           | 🗇 Ground      | lwater l         | 🗆 Surfac           | e Wa          | ater         |             | ∕<br>Othe       | er S.o        | n+i                 | . L                   | later                                  | 1           | Date      | 24/97                                 |
| Sampling Specifications       Initial Measurements         Requested Sampling  | roject Nam            | ne h            | JES           | Mila             |                    | -             | +;,          | . Ţ         | -10             | (1)           | <i>p_</i>           | <u> </u>              | P                                      | -<br>roiect | No. 1     | 7309                                  |
| Sampling Specifications       Initial Measurements         Requested Sampling  | roiect Man            | nager           | CM            | Clark            | gr v               | - <del></del> | <i></i>      |             | _ <u>`</u> _\   | -\ <b>f</b> / |                     |                       | P                                      | hase.       | Task No   | 6000 77                               |
| Sampling Specifications       Initial Measurements         Requested Sampling  | ite Name              | N               | Ailay         | ro Pl            | unt,               | BI            | DDw          | Fiel        | D,              | NN            | 1                   |                       | ······································ |             |           |                                       |
| Requested Wait Following       Initial Water Depth (feet)         Development/Purging (hours)  | Sampling<br>Requested | Speci<br>Sampli | fication:     | S                |                    |               | Ini          | tial I      | Me              | asuren        | nen                 | ts                    |  |             |           |                                       |
| Nater Quality/Water Collection       D0 = Dissolved Oxygen; Cond = Cond         Date       Time       Sampler<br>Initials       Water Quality Readings       Water Collection Data       Note<br>Explain in Collection         Date       Time       Final       Final       Final       Final         Temp, pH       D0       Cond.       Removed<br>(galions)       Removed<br>(galions)       Final       Ball<br>Water         Container       Temp, Final       Cond.       Galianes)       Galianes       Final       Depth fill         Container       Type:       G = Clear Glass; A = Amber Glass; F = Plastic; V = VOA Vial (Glass); 0 = Other (Specify); = None         Container       Preservatives: H = HO! N = HNO; S = H:SO; A = NaOH: 0 = Other (Specify); = None         Parameter List       Container       Field       Preserved       Colled       During       Contents         Preserved       Colled       During       Colled       During       Container         Primeter List       Number       Type       Volume (nt.1)       Yes       No       Sample 1.0.       Time, Final         Metals       J       P       SSO       V       HMD; V       WES-ST-D:D:D:D       Second 1.0.       Second 1.0.         VD A'3       Z       YO       YO   | Requested             | Wait F          | ollowing      |                  |                    |               |              |             |                 |               |                     |                       |  |             |           | · · · · · · · · · · · · · · · · · · · |
| Date     Sampler<br>Initials     Water Quality Readings     Water Collection Data     Initials<br>(Explain in C<br>Below       Temp.     pH     D0     Cond.     Permoval     Pump     Bail     Water     Depth (tt)       Cond.     Temp.     pH     D0     Cond.     (galfmin)     Depth (tt)     Bail     Water       Container     Temp.     pH     D0     Cond.     (galfmin)     Depth (tt)     Depth (tt)       Container Type:     G = Clear Glast, A = Amber Glast;     P = Plastic;     V = VOA Vial (Glast);     0 = Other (Specify);     = None       Analytical<br>Parameter List     Container     Field     Preserved     Cooled     Content       Number     Type     Volume (nL1)     Yes     No     WES-ST-/2497     100/s       VD A <sup>1</sup> 3     V     YD     HCL     V     V       Filter Type     MA     IDDU  |                       |                 |               |                  |                    | -             |              | Nona        | que             | eous Lia      | luids               | Pres                  |  |             |           | ond - Conductivi                      |
| Date       Time       Initials       Water Ouality Readings       Water Collection Data       Initials       Init   |                       |                 |               | 1                | <u> </u>           |               |              |             |                 |               |                     |                       |  |             | oxygen, c | Notes                                 |
| Temp.       pH       DO       Cond.       Volume       Removal       Pump       Bail       Final         Col       (mg(L)       (multication)       (galfone)       (galfone)       Dopth (ft)       Depth (ft)         Container       Container       (galfone)       (galfone)       (galfone)       Depth (ft)         Container       Type:       G = Clear Glass;       A = Amber Glass;       P = Plastic;       V = VOA Vial (Glass);       O = Other (Specify);         Sample Containers       Preservatives:       H = HCI:       N = HNO;       S = HJSO4;       A = NaOH;       O = Other (Specify);       = None         Analytical       Container       Field       Preserved       Coolection       Contents         Net al S       J       P       2SO       V       HMD;       W/SS-ST-//, 9/97       /////, 0/1,         VD A'1       3       V       4D       V       H(L)       V       V       J         Senje - J/DA'1       A       IDDD       -       V       J<   | Date                  | _ Time          |               |                  | ater Qua           | lity R        | eadi         | ngs         |                 |               | Wa                  | ter C                 | ollection                              | Data        |           | (Explain in Comme<br>Below)           |
| Container Type: G = Clear Glass; A = Amber Glass; P = Plastic; V = VOA Vial (Glass); O = Other (Spectry I; = None         Container Type: G = Clear Glass; A = Amber Glass; P = Plastic; V = VOA Vial (Glass); O = Other (Spectry I; = None         Analytical Parameter List       Container         Number       Type         VDA'S       3         VDA'S       3         VDA'S       3         Analytical Parameter List       Passo         VDA'S       3         VDA'S       4         IDDD   |                       |                 | +             | $\sim$           | ρН                 |               |              |             | · 1             | Removed       | R                   | ate                   | Intake                                 | Bail        | Water     |                                       |
| Sample Containers       Preservatives: H = HCI: N = HNO2; S = H/SO4; A = NaOH; O = Other (Specify I; = None         Analytical Parameter List       Container       Field Filtered       Preserved Collection       Cooled During Collection       Comments         M etals       1       P       2SD       V       HMD2       WES-ST-D497       100D         VD A <sup>13</sup> 3       V       4D       V       HCL       V       V         Semj-UDA <sup>15</sup> 2       A       1DDD       -       V       V       V         Filter Type       MA       DDD       -       V       Container       Semple Container       Semple Container       Semple Container       Time   |                       |                 |               |                  |                    | - ungr        |              |             | cini            | (gailons)     | iyai                | /110(1)               | Deptir (it)                            |             | Depurity  |                                       |
| Sample Containers       Preservatives: H = HCI: N = HNO2; S = H2SO4; A = NaOH; O = Other (Specify I; = None         Analytical<br>Parameter List       Container       Field<br>Filtered       Preserved<br>Filter Type       Cooled<br>During<br>Collection       Comments         M et al S       1       P       2SD       V       HMD2       WESS-ST-D497       100D         VD A <sup>13</sup> 3       V       4D       V       HCL       V       V         Semj-UDA <sup>15</sup> 2       A       1DDD       -       V       V       V         Filter Type       MA       Container       Container       Container       Container       Container       Container       Container       Container         M et al S       1       P       2SD       V       HMD2       WESS-ST-D497       100D         VD A <sup>15</sup> 3       V       4D       V       HCL       V       V       V         Semj-UDA <sup>15</sup> 2       A       1DDD       -       V       V       V       V         Filter Type       MA       Chain-of-Custody Form Number       C3D2D       C3D2D   |                       |                 |               |                  | $\left  - \right $ |               | 4            | $\vee \neq$ | $\mathcal{H}$   |               |                     |                       |  |             |           |                                       |
| Cample Containers       Preservatives: H = HCI: N = HNO; S = H;SO4: A = NaOH; O = Other (Specify ); = None         Analytical<br>Parameter List       Container       FileId<br>Filtered       Preserved       Cooled<br>During<br>Collection       Comments         Metals       1       P       2SD       V       HMD;       WES-ST-D:9497       100D         VDA's       3       V       4D       V       H(L       V       V         Semi-WA's       2       A       10DD       -       V       V       V         Filter Type       MA       Chain-of-Custody Form Number       C3D2D       Chain-of-Custody Form Number       C3D2D   |                       |                 |               |                  |                    |               |              |             | $ \rightarrow $ | ~             |                     |                       |  | ļ           |           |                                       |
| Comple Containers       Preservatives: H = HCI: N = HNO2: S = H2O4: A = NaOH: O = Other (Specify ): = None         Analytical<br>Parameter List       Container       Filed<br>Filed       Preserved       Cooled<br>During<br>Collection       Comments         M et al S       1       P       2SD       V       HMD2       WES-ST-D497       100D         VD A <sup>13</sup> 3       V       4D       V       HCL       V       V         Semi-UDA <sup>15</sup> 2       A       10DD       -       V       V       V         Filter Type       MA       Container       Container       Container       Container       Container       Container       Container         M et al S       1       P       2SD       V       HMD2       WES-ST-D497       100D         VD A <sup>15</sup> 3       V       4D       V       HCL       V       V         Semi-UDA <sup>15</sup> A       10DD       -       V       V       V       V         Filter Type       MA       Chain-of-Custody Form Number       C3D2D       C3D2D       Chain-of-Custody Form Number       C3D2D   |                       |                 |               |                  |                    |               |              |             |                 |               | $\downarrow$        |                       |  |             |           |                                       |
| Sample Containers       Preservatives: H = HCI: N = HNO2; S = H2SO4; A = NaOH; O = Other (Specify I; = None         Analytical<br>Parameter List       Container       Field<br>Filtered       Preserved<br>Filter Type       Cooled<br>During<br>Collection       Comments         M et al S       1       P       2SD       V       HMD2       WESS-ST-D497       100D         VD A <sup>13</sup> 3       V       4D       V       HCL       V       V         Semj-UDA <sup>15</sup> 2       A       1DDD       -       V       V       V         Filter Type       MA       Container       Container       Container       Container       Container       Container       Container       Container         M et al S       1       P       2SD       V       HMD2       WESS-ST-D497       100D         VD A <sup>15</sup> 3       V       4D       V       HCL       V       V       V         Semj-UDA <sup>15</sup> 2       A       1DDD       -       V       V       V       V         Filter Type       MA       Chain-of-Custody Form Number       C3D2D       C3D2D   |                       | ·               |               |                  |                    |               |              |             |                 |               |                     |                       |  |             |           |                                       |
| Sample Containers       Preservatives: H = HCI: N = HNO2: S = H2SO4: A = NaOH; O = Other (Specify I; = None         Analytical<br>Parameter List       Container       Field<br>Filtered       Preserved<br>Filter Type       Cooled<br>During<br>Collection       Comments         M et al S       1       P       2SD       V       HMD2       WES-ST-D497       100D         VD A <sup>13</sup> 3       V       4D       V       HCL       V       V         Semi-UDA <sup>15</sup> 2       A       1DDD       -       V       V       V         Filter Type       MA       Container       Container       Container       Container       Container       Container       Container       Container         M et al S       1       P       2SD       V       HMD2       WES-ST-D497       100D         VD A <sup>15</sup> 3       V       4D       V       HCL       V       V       V         Semi-UDA <sup>15</sup> 2       A       1DDD       -       V       V       V       V       V         Filter Type       MA       Chain-of-Custody Form Number       C3D2D       C3D2D       Chain-of-Custody Form Number       C3D2D  |                       |                 |               |                  |                    |               |              |             |                 |               |                     |                       |  |             |           |                                       |
| VI et al s 1 P 2SD V HND2 WES-ST-12497 1000<br>VDA'S 3 V 40 HCL V V<br>Semi-WDA'S 2 A IDDU - V V V<br>Semi-WDA'S 2 A IDDU - V V V<br>Chain-of-Custody Form Number 302D   | Analytic              | al              |               | Containe         | er                 |               | Fie<br>Filte | eld<br>ered |                 |               | Coo<br>Dur<br>Colle | oled<br>ing<br>ction_ |  |             | Comme     | nts                                   |
| VDA's     3     V     40     HCL       Semi-UDA's     2     A     IDDU   Filter Type NA Chain-of-Custody Form Number 3D2.D   | IN/ I                 |                 | Number        | Туре             |                    |               | Yes          | No<br>/     | 11              | . 1.          | Yes                 | No                    | Sampla                                 | <u>1.0</u>  |           | Timp                                  |
| Seni-UDA <sup>S</sup> Q A IDDU - V<br>Seni-UDA <sup>S</sup> Q A IDDU - V<br>Seni-UDA <sup>S</sup> Q A IDDU - V<br>Seni-VDA <sup>S</sup> Q A ID | Metals                | s               |               | <u> </u>         | asi                | 2             |              |             | <u>  /]</u>     | ND2           |                     | ·                     | WES-S                                  | <u>T-1</u>  | 497       | 1000                                  |
| Filter Type  | VDA'                  |                 |               | V                | 40                 | <u> </u>      |              | V           | <i>†</i>        | <u>I</u> (L   |                     |                       | ļ                                      | 1           |           |                                       |
|  | Semi-VL               | JA'S            | 2             | A                | IDDL               | )             |              | /           |                 | -             | V                   |                       |  | V_          |           | $\downarrow$                          |
|  |                       |                 | l             |                  |                    |               |              |             |                 | 1             |                     |                       |  |             |           |                                       |
|  | ,,,,,,,,,,,           |                 |               |                  |                    |               |              |             |                 |               |                     |                       |  |             |           |                                       |
|  |                       |                 |               |                  |                    |               | c            |             |                 |               |                     |                       |  |             |           |                                       |
|  |                       |                 |               |                  |                    | _             |              | K           | Z               | 5.1           |                     |                       |  |             | <u>.</u>  |                                       |
|  |                       |                 |               |                  |                    |               |              |             |                 |               | 2                   |                       | <u></u>                                |             | /*        |                                       |
|  |                       |                 |               |                  |                    |               |              | }           |                 |               |                     |                       |  |             |           |                                       |
|  |                       |                 |               |                  |                    |               |              |             |                 |               |                     |                       |  |             |           |                                       |
|  | Filter Type           |                 | NA            |                  |                    |               |              |             | C               | Chain-of      | -Cus                | todv                  | Form Ni                                | ımber       | C         | 5020                                  |
| Comments   | _                     |                 | •             |                  |                    |               |              |             | - '             |               |                     |                       |  |             |           |                                       |
|  | Commen                | ts              |               |                  | ·······            |               |              |             |                 |               |                     | ·                     | <u></u>                                |             |           |                                       |
|  |                       |                 |               |                  |                    |               |              |             |                 |               |                     |                       |  |             |           |                                       |
|  |                       |                 |               |                  |                    |               |              |             |                 |               |                     |                       |  |             |           |                                       |
|  |                       | $\sim$          |               | $\frown$ $\land$ |                    |               |              |             |                 |               | Τ.                  |                       |  |             |           |                                       |
| Signature Date 1/24/97 Reviewer Date   | Signature_            |                 | sm.           | Cha              | ny                 |               |              |             | _ [             | Date_]        | 12                  | 41                    | 97 R                                   | eview       | /er       | Date                                  |
|  |                       |                 | $^{-}\Lambda$ |                  | •                  |               |              |             |                 |               |                     |                       |  |             |           |                                       |

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**APPENDIX B - LABORATORY REPORTS** 

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5555 North Service Road Burlington, Ontario, Canada L7L 5H7 Tel: (905) 332-8788 Fax: (905) 332-9169

# Certificate of Analysis

### **CLIENT INFORMATION**

#### LABORATORY INFORMATION

| Attention:           | Cory Chance               | Contact:        | Ada Blythe, B.Sc., C.Chem. |
|----------------------|---------------------------|-----------------|----------------------------|
| <b>Client Name:</b>  | Philip Environmental Inc. | Project:        | AN961368                   |
| Project:             | 17039                     | Date Received:  | 96/12/13                   |
| <b>Project Desc:</b> | WFS Septic Leach Field    | Date Reported:  | 97/01/02                   |
| Address:             | 4000 Monroe Road          | Submission No.: | 6L0416                     |
|                      | Farmington, NM            | Sample No.:     | 055066-055069              |
|                      | 87401                     |                 |                            |
| Fax Number:          | 505 326-2388              |                 |                            |
| Phone Number:        | 505 326-2262              |                 |                            |

**NOTES:** '-' = not analysed '<' = less than Method Detection Limit (MDL) 'NA' = no data available LOQ can by determined for all analytes by multiplying the appropriate MDL X 3.33 Solids data is based on dry weight except for biota analyses. Organic analyses are not corrected for extraction recovery standards except for isotope dilution methods, (i.e. CARB 429 PAH, all PCDD/F and DBD/DBF analyses)

Methods used by Zenon are based upon those found in 'Standard Methods for the Examination of Water and Wastewater', Seventeenth Edition. Other methods are based on the principles of MISA or EPA methodologies.

All work recorded herein has been done in accordance with normal professional standards using accepted testing methodologies, quality assurance and quality control procedures except where otherwise agreed to by the client and testing company in writing. Any and all use of these test results shall be limited to the actual cost of the pertinent analysis done. There is no other warranty expressed or implied. Your samples will be retained at Zenon for a period of three weeks from receipt of data or as per contract.

COMMENTS:

Certified by:

Page 1

A division of PHILIP Analytical Services Corporation

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|                                 | Client 1D:<br>Zenon 1D:<br>Date Sampled: |       | Method<br>Blank<br>055065 96<br>96/12/12 | BH1-3-5<br>055066 96<br>96/12/12 | BH1<br>10-12<br>055067 96<br>96/12/12 | BH2<br>15-17<br>055068 96<br>96/12/12 |  |
|---------------------------------|--|-------|--|----------------------------------|---------------------------------------|---------------------------------------|--|
| Component                       | MDL                                      | Units |  |                                  |                                       |                                       |  |
| Mercury via SW846 Method 7471   |  |       |  |                                  |                                       |                                       |  |
| Mercury                         | 0.04                                     | mg/kg | <  | <                                | <                                     | <                                     |  |
| Metals via SW846 Method 6010    |  |       |  |                                  |                                       |                                       |  |
| Arsenie                         | 5.0                                      | mg/kg | <10                                      | <10                              | <10                                   | <10                                   |  |
| Selenium                        | 10                                       | "     | <20                                      | <20                              | <20                                   | <20                                   |  |
| Barium                          | 0.1                                      | u     | <0.2                                     | 280                              | 130                                   | 190                                   |  |
| Cadmium                         | 0.2                                      | "     | <0.4                                     | <0.4                             | <0.4                                  | <0.4                                  |  |
| Chromium                        | 5  | 11    | <10                                      | <10                              | <10                                   | <10                                   |  |
| Lead                            | 10                                       | н     | <20                                      | <20                              | <20                                   | <20                                   |  |
| Silver                          | 0.5                                      | "     | <1.0                                     | <1.0                             | <1.0                                  | <1.0                                  |  |
| Volatiles via SW846 Method 8260 |  |       |  |                                  |                                       |                                       |  |
| Acetone                         | 0.030                                    | mg/kg | 0.17                                     | 0.18                             | 0.15                                  | 0.14                                  |  |
| Aerolein                        | 0.010                                    | "     | <  | <                                | <                                     | <                                     |  |
| Acrylonitrile                   | 0.010                                    | н     | <  | <                                | <                                     | <                                     |  |
| Benzene                         | 0.005                                    | n     | <  | <                                | <                                     | <                                     |  |
| Bromoform                       | 0.010                                    | "     | <  | <                                | <                                     | <                                     |  |
| Bromomethane                    | 0.010                                    | н     | <  | <                                | <                                     | <                                     |  |
| 2-Butanone                      | 0.015                                    |       | <  | <                                | <                                     | <                                     |  |
| Carbon Disulfide                | 0.010                                    | "     | <  | <                                | <                                     | <                                     |  |
| Carbon Tetrachloride            | 0.010                                    | 0     | <  | <                                | <                                     | <                                     |  |
| Chlorobenzene                   | 0.005                                    | "     | <  | <                                | <                                     | <                                     |  |
| Chlorodibromomethane            | 0.005                                    | ч     | <  | <                                | <                                     | <                                     |  |
| Chloroethane                    | 0.010                                    | "     | <  | <                                | <                                     | <                                     |  |
| 2-Chloroethylvinylether         | 0.010                                    | 11    | <  | <                                | <                                     | <                                     |  |
| Chloroform                      | 0.005                                    | 55    | <  | <                                | <                                     | <                                     |  |
| Chloromethane                   | 0.010                                    | 4     | <  | <                                | <                                     | <                                     |  |
| 1,2-Dichlorobenzene             | 0.005                                    | b.    | <  | <                                | <                                     | <                                     |  |
| 1,3-Dichlorobenzene             | 0.005                                    | "     | <  | <                                | <                                     | <                                     |  |
| 1,4-Dichlorobenzene             | 0.005                                    | "     | <  | 0.005                            | <                                     | <                                     |  |
| Dichlorobromomethane            | 0.005                                    | н     | <  | <                                | <                                     | <                                     |  |
| 1,1-Dichloroethane              | 0.005                                    | T)    | <  | <                                | <                                     | <                                     |  |
| 1,2-Dichloroethane              | 0.005                                    | **    | <  | <                                | . <                                   | <                                     |  |
| 1,1-Dichloroethene              | 0.010                                    | **    | <  | <                                | <                                     | <                                     |  |
| cis-1,2-Dichloroethene          | 0.010                                    | 0     | <  | <                                | <                                     | <                                     |  |
| trans-1,2-Dichloroethene        | 0.010                                    | 15    | <  | <                                | <                                     | <                                     |  |
| 1,2-Dichloropropane             | 0.005                                    | **    | <  | <                                | <                                     | <                                     |  |
| cis-1,3-Dichloropropene         | 0.005                                    | 0     | <  | <                                | <                                     | <                                     |  |
| trans-1,3-Dichloropropene       | 0.005                                    | в     | <  | <                                | <                                     | <                                     |  |
| Ethylbenzene                    | 0.005                                    | ۴r    | <  | <                                | <                                     | <                                     |  |
| 2-Hexanone                      | 0.010                                    | н     | <  | <                                | <                                     | < .                                   |  |
| Methylene Chloride              | 0.010                                    | "     | 0.020                                    | 0.039                            | 0.034                                 | 0.029                                 |  |
| 4-Methyl-2-Pentanone            | 0.010                                    | u     | <  | <                                | <                                     | <                                     |  |
| Styrene                         | 0.005                                    |       | <  | <                                | <                                     | <                                     |  |
| 1,1,1,2-Tetrachloroethane       | 0.010                                    | и     | <  | <                                | <                                     | ~                                     |  |
| 1,1,2,2-Tetrachloroethane       | 0.010                                    | н     | <  | <                                | <                                     | <                                     |  |
| Tetrachloroethene               | 0.020                                    | 11    | <  | < .                              | <                                     | <                                     |  |
| i oquonitor oculono             | 0.040                                    |       | · ·                                      | ~                                | · ·                                   | -                                     |  |

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|                                    | Client ID:<br>Zenon ID:<br>Date Sampled: |       | Method<br>Blank<br>055065 96<br>96/12/12 | BH1-3-5<br>055066 96<br>96/12/12        | BH1<br>10-12<br>055067 96<br>96/12/12 | BH2<br>15-17<br>055068 96<br>96/12/12 |
|------------------------------------|--|-------|--|---|---------------------------------------|---------------------------------------|
| Component                          | MDL                                      | Units |  |   |                                       |                                       |
| 1,1,1-Trichloroethane              | 0.005                                    | 11    | <  | <                                       | <                                     | <                                     |
| 1,1,2-Trichloroethane              | 0.010                                    |       | <  | <                                       | <                                     | <                                     |
| Trichloroethene                    | 0.005                                    | U.    | <  | <                                       | <                                     | <                                     |
| Vinyl Acetate                      | 0.010                                    | 11    | <  | <                                       | <                                     | <                                     |
| Vinyl Chloride                     | 0.010                                    | ц     | <  | <                                       | <                                     | <                                     |
| Xylenes(Total)                     | 0.005                                    | "     | <  | <                                       | <                                     | <                                     |
| Surrogate Recoveries               |  | %     |  |   |                                       |                                       |
| d4-1,2-Dichloroethane              |  |       | 95                                       | 97                                      | 103                                   | 98                                    |
| d8-Toluene                         |  |       | 98                                       | 99                                      | 94                                    | 98                                    |
| Bromofluorobenzene                 |  |       | 95                                       | 95                                      | 95                                    | 91                                    |
| Semi Volatiles via SW846 Method 82 | 27()                                     |       | ,,,                                      | ,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,, |                                       |                                       |
| Phenol                             | 0.11                                     | mg/kg | < 0.22                                   | < 0.22                                  | < 0.22                                | <0.22                                 |
| Bis(2-chloroethyl)ether            | 0.18                                     | "     | < 0.36                                   | < 0.36                                  | < 0.36                                | < 0.36                                |
| 2-Chlorophenol                     | 0.27                                     | n.    | < 0.54                                   | < 0.54                                  | <0.54                                 | < 0.54                                |
| 1,3-Dichlorobenzene                | 0.20                                     |       | <().4()                                  | <0.40                                   | < 0.40                                | < 0.40                                |
| 1-4-Dichlorobenzene                | 0.20                                     |       | <0.40                                    | < 0.40                                  | < 0.40                                | <0.40                                 |
| 1,2-Dichlorobenzene                | 0.20                                     | 0     | < 0.40                                   | <0.40                                   | < 0.40                                | < 0.40                                |
| Bis(2-chloroisopropyl)ether        | 0.15                                     |       | < 0.30                                   | < 0.30                                  | < 0.30                                | < 0.30                                |
| Hexachloroethane                   | 0.20                                     | н     | < 0.40                                   | <0.40                                   | <0.40                                 | < 0.40                                |
| N-Nitroso-di-N-Propylamine         | 0.21                                     |       | < 0.42                                   | < 0.42                                  | < 0.42                                | < 0.42                                |
| Nitrobenzene                       | 0.20                                     |       | <0.40                                    | < 0.40                                  | <0.40                                 | <0.40                                 |
| Isophorone                         | 0,40                                     | •     | < 0.80                                   | < 0.80                                  | < 0.80                                | <0.80                                 |
| 2-Nitrophenol                      | 0.14                                     | "     | <0.28                                    | < 0.28                                  | <0.28                                 | <0.28                                 |
| 2,4-Dimethylphenol                 | 0.17                                     | р     | <0.34                                    | < 0.34                                  | < 0.34                                | < 0.34                                |
| Bis(2-chloroethoxy)methane         | 0.13                                     | "     | < 0.26                                   | < 0.26                                  | < 0.26                                | <0.26                                 |
| 2,4-Dichlorophenol                 | 0.12                                     | ų     | <0.20                                    | <0.24                                   | <0.24                                 | <0.24                                 |
| 1,2,4-Trichlorobenzene             | 0.12                                     | 11    | <0.24                                    | <0.40                                   | <0.40                                 | <0.40                                 |
| Naphthalene                        | 0.03                                     |       | <0.46                                    | <0.46                                   | <0.46                                 | <0.06                                 |
| Hexachlorobutadiene                | . 0.20                                   |       | <0.40                                    | <0.40                                   | <0.40                                 | <0.40                                 |
| 4-Chloro-3-Methylphenol            | 0.14                                     | ۶t    | <0.40                                    | <0.40                                   | <0.40                                 | <0.48                                 |
| Hexachlorocyclopentadiene          | 0.14                                     | 0     | <0.28                                    | <0.28                                   | <0.28                                 | <0.28                                 |
| 2,4,6-Trichlorophenol              | 0.12                                     | "     | <0.40                                    | <0.40                                   | <0.40                                 | <0.40                                 |
| -                                  | 0.12                                     | н     | <0.24                                    | <0.24                                   | <0.24                                 | <0.24                                 |
| 2-Chloronaphthalene                | 0.09                                     |       | <0.18                                    | <0.18                                   | <0.18                                 | <0.18                                 |
| Acenaphthylene                     |  | **    |  |   | <0.08                                 |                                       |
| Dimethyl phthalate                 | 0.11                                     | u.    | <0.22                                    | < 0.22                                  | <0.22                                 | <0.22<br><0.12                        |
| 2,6-Dinitrotoluene                 | 0.06<br>0.07                             |       | <0.12<br><0.14                           | <0.12<br><0.14                          | <0.12                                 | <0.12                                 |
| Acenaphthene<br>2,4-Dinitrophenol  |  | **    |  |   | <0.14<br><0.96                        | <0.14                                 |
|                                    | 0.48                                     | u     | <0.96                                    | <0.96                                   |                                       |                                       |
| 2,4-Dinitrotoluene                 | 0.05                                     |       | <0.10                                    | <0.10                                   | <0.10                                 | <0.10                                 |
| 4-Nitrophenol                      | 0.14                                     | F1    | <0.28                                    | <0.28                                   | <0.28                                 | <0.28                                 |
| Fluorene                           | 0.03                                     |       | <0.06                                    | <0.06                                   | <0.06                                 | < 0.06                                |
| 4-Chlorophenylphenylether          | 0.09                                     | 59    | <0.18                                    | <0.18                                   | <0.18                                 | <0.18                                 |
| Diethyl phthalate                  | 0.11                                     |       | <0.22                                    | <0.22                                   | <0.22                                 | <0.22                                 |
| 4,6-Dinitro-2-methylphenol         | 0.15                                     | **    | < 0.30                                   | < 0.30                                  | < 0.30                                | <0.30                                 |
| N-Nitrosodiphenylamine             | 0.19                                     |       | <0.38                                    | < 0.38                                  | <0.38                                 | < 0.38                                |
| 4-Bromophenylphenylether           | 0.03                                     | "     | <0.06                                    | <0.06                                   | <0.06                                 | <0.06                                 |
| Hexachlorobenzene                  | 0.20                                     |       | <0.40                                    | <0.40                                   | <0.40                                 | <().4()                               |

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## Zenon Environmental Laboratories - Certificate of Analysis

|                            | Client ID:<br>Zenon ID: |            | Method<br>Blank<br>055065 96  | BH1-3-5<br>055066 96 | BH1<br>10-12<br>055067 96 | BH2<br>15-17<br>055068 96 |
|----------------------------|-------------------------|------------|---|----------------------|---------------------------|---------------------------|
|                            | Date Sampled:           |            | 96/12/12  | 96/12/12             | 96/12/12                  | 96/12/12                  |
| Component                  | MDL                     | Units      |   |                      |                           |                           |
| Pentachlorophenol          | 0.11                    | 11         | < 0.22  | < 0.22               | < 0.22                    | <0.22                     |
| Phenanthrene               | 0.03                    | п          | < 0.06  | < 0.06               | <0.06                     | < 0.06                    |
| Anthracene                 | 0.02                    |            | < 0.04  | < 0.04               | < 0.04                    | <0.04                     |
| Di-n-butyl phthalate       | 0.11                    | н          | < 0.22  | < 0.22               | < 0.22                    | < 0.22                    |
| Fluoranthene               | 0.02                    | н          | < 0.04  | < 0.04               | < 0.04                    | < 0.04                    |
| Pyrene                     | 0.03                    | "          | < 0.06  | < 0.06               | <0.06                     | <0.06                     |
| Benzyl butyl phthalate     | 0.06                    | **         | < 0.12  | < 0.12               | <0.12                     | < 0.12                    |
| 3,3-Dichlorobenzidine      | 0.10                    | U          | <0.20   | < 0.20               | < 0.20                    | <0.20                     |
| Benzo(a)anthracene         | 0.02                    | •          | < 0.04  | < 0.04               | < 0.04                    | < 0.04                    |
| Chrysene                   | 0.03                    | **         | < 0.06  | < 0.06               | < 0.06                    | <0.06                     |
| Bis(2-ethylhexyl)phthalate | 0.14                    |            | < 0.28  | < 0.28               | < 0.28                    | <0.28                     |
| Di-n-octyl phthalate       | 0.11                    | "          | < 0.22  | < 0.22               | < 0.22                    | < 0.22                    |
| Benzo(b)fluoranthene       | 0.04                    | н          | .<0.08  | <0.08                | < 0.08                    | <0.08                     |
| Benzo(k)fluoranthene       | 0.04                    |            | < 0.08  | < 0.08               | < 0.08                    | < 0.08                    |
| Benzo(a)pvrene             | 0.05                    | "          | < 0.10  | <0.10                | <0.10                     | <0.10                     |
| Indeno(1,2,3-cd)pyrene     | 0.06                    | ц          | < 0.12  | < 0.12               | <0.12                     | < 0.12                    |
| Dibenzo(a,h)anthracene     | 0.04                    | "          | < 0.08  | < 0.08               | < 0.08                    | < 0.08                    |
| Benzo(ghi)perylene         | 0.04                    | **         | < 0.08  | < 0.08               | < 0.08                    | < 0.08                    |
| N-Nitrosodimethylamine     | 1.0                     | mg/kg      | <2.0  | <2.0                 | <2.0                      | <2.0                      |
| Aniline                    | 0.50                    | "          | <1.0  | <1.0                 | <1.0                      | <1.0                      |
| Benzyl alcohol             | 0.50                    | и          | <1.0  | <1.0                 | <1.0                      | <1.0                      |
| Carbazole                  | 0.50                    | Ħ          | <1.0  | <1.0                 | <1.0                      | <1.0                      |
| 2-Methylphenol             | 0.50                    | II.        | <1.0  | <1.0                 | <1.0                      | <1.0                      |
| Benzoic acid               | 0.50                    | 15         | <i.0< td=""><td>&lt;].0</td><td>&lt;1.0</td><td>&lt;1.0</td></i.0<> | <].0                 | <1.0                      | <1.0                      |
| 4-Chloroaniline            | 0.50                    | ¢¥         | <1.0  | <1.0                 | <1.0                      | <1.0                      |
| 2-Methylnaphthalene        | 0.10                    | <b>1</b> 1 | < 0.20  | <0.20                | <0.20                     | < 0.20                    |
| 2,4,5-Trichlorophenol      | 0.10                    | н          | <0.20   | <0.20                | <0.20                     | <0.20                     |
| 2-Nitroaniline             | 0.50                    | <b></b>    | <1.0  | <1.0                 | <1.0                      | <1.0                      |
| 3-Nitroaniline             | 0.50                    | "          | <1.0  | 0.1>                 | <1.0                      | <1.0                      |
| Dibenzofuran               | 0.50                    | н          | <1.0  | <1.0                 | <1.0                      | <1.0                      |
| Benzidine                  | 0.50                    | н          | <1.0  | <1.0                 | <1.0                      | <1.()                     |
| 4-Nitroaniline             | 0.50                    | f*         | <1.0  | <1.0                 | <1.()                     | <1.0                      |
| Surrogate Recoveries       |                         | %          |   |                      |                           |                           |
| 2-Fluorophenol             |                         |            | 25  | 40                   | 39                        | 43                        |
| d5-Phenol                  |                         |            | 58  | 68                   | 66                        | 65                        |
| d5-Nitrobenzene            |                         |            | 19  | 42                   | 41                        | 43                        |
| 2-Fluorobiphenyl           |                         |            | 73  | 59                   | 60                        | 62                        |
| 2.4,6-Tribromophenol       |                         |            | 101   | 69                   | 61                        | 64                        |
| d14-p-Terphenyl            |                         |            | 105   | 75                   | 72                        | 70                        |

| Component                       | Client ID:<br>Zenon ID:<br>Date Sampled:<br>MDL | Units | BH3<br>13-15<br>055069 96<br>96/12/12 | BH3<br>13-15<br>055069 96<br>96/12/12<br>MS | BH3<br>13-15<br>055069 96<br>96/12/12<br>% recoveries | BH3<br>13-15<br>055069 96<br>96/12/12<br>MSD |
|---------------------------------|---|-------|---------------------------------------|---|---|--|
| Mercury via SW846 Method 7471   |   |       |                                       |   |   |  |
| Mercury                         | 0.04  | mg/kg | <                                     | -   | -   | -  |
| Metals via SW846 Method 6010    |   |       |                                       |   |   |  |
| Arsenic                         | 5.0   | mg/kg | <10                                   | -   | -   | -  |
| Selenium                        | 10  | 11    | <20                                   | -   | -   | -  |
| Barium                          | 0.1   | н     | 1501                                  | -   | -   | -  |
| Cadmium                         | 0.2   | "     | <()_4                                 | -   | -   | ~  |
| Chromium                        | 5   | 11    | <10                                   | -   | -   | -  |
| Lead                            | 10  | **    | <20                                   | -   | -   | -  |
| Silver                          | 0.5   | н     | <1.0                                  | -   | -   | -  |
| Volatiles via SW846 Method 8260 |   |       |                                       |   |   |  |
| Acetone                         | 0.030   | mg/kg | 0.12                                  | 0.4 I                                       | 160   | 0.33   |
| Acrolein                        | 0.010   | n     | <                                     | 0.42  | 83  | 0.36   |
| Acrylonitrile                   | 0.010   | "     | <                                     | 0.40  | 80  | 0.39   |
| Benzene                         | 0.005   | п     | <                                     | 0.23  | 94  | 0.23   |
| Bromoform                       | 0.010   | **    | <                                     | 0.26  | 110   | 0.26   |
| Bromomethane                    | 0.010   | u     | <                                     | 0.29  | 120   | 0.29   |
| 2-Butanone                      | 0.015   | "     | <                                     | 0.21  | 83  | 0.19   |
| Carbon Disulfide                | 0.010   | 0     | <                                     | 0.22  | 88  | 0.20   |
| Carbon Tetrachloride            | 0.010   |       | <                                     | 0.26  | 110   | 0.26   |
| Chlorobenzene                   | 0.005   | ()    | <                                     | 0.23  | 91  | 0.23   |
| Chlorodibromomethane            | 0.005   |       | <                                     | 0.25  | 100   | 0.25   |
| Chloroethane                    | 0.010   | 9     | <                                     | 0.27  | 110   | 0.25   |
| 2-Chloroethylvinylether         | 0.010   | "     | <                                     | 0.16  | 64  | 0.15   |
| Chloroform                      | 0.005   | ч     | <                                     | 0.24  | 95  | 0.23   |
| Chloromethane                   | 0.010   | "     | <                                     | 0.27  | 110   | 0.24   |
| 1,2-Dichlorobenzene             | 0:005   | н     | <                                     | 0.24  | 94  | 0.23   |
| 1,3-Dichlorobenzene             | 0.005   | 11    | <                                     | 0.22  | 88  | 0.21   |
| 1,4-Dichlorobenzene             | 0.005   | н     | <                                     | 0.22  | 87  | 0.22   |
| Dichlorobromomethane            | 0.005   |       | <                                     | 0.25  | 100   | 0.25   |
| 1,1-Dichloroethane              | 0.005   | п     | <                                     | 0.23  | 93  | 0.22   |
| 1,2-Dichloroethane              | 0.005   | н     | <                                     | 0.24  | 97  | 0.23   |
| 1,1-Dichloroethene              | 0.010   | "     | <                                     | 0.24  | 97  | 0.22   |
| cis-1,2-Dichloroethene          | 0.010   | п     | <                                     | 0.22  | 89  | 0.21   |
| trans-1,2-Dichloroethene        | 0.010   | и     | <                                     | 0.22  | 88  | 0.21   |
| 1,2-Dichloropropane             | 0.005   | "     | <                                     | 0.23  | 92  | 0.24   |
| cis-1,3-Dichloropropene         | 0.005   | 11    | <                                     | 0.23  | 90  | 0.23   |
| trans-1,3-Dichloropropene       | 0.005   | "     | <                                     | 0.22  | 89  | 0.23   |
| Ethylbenzene                    | 0.005   | п     | <                                     | 0.23  | 91  | 0.23   |
| 2-Hexanone                      | 0.010   |       | <                                     | 0.18  | 73  | 0.18   |
| Methylene Chloride              | 0.010   | н     | 0.033                                 | 0.25  | 99  | 0.24   |
| 4-Methyl-2-Pentanone            | 0.010   | **    | <                                     | 0.20  | 80  | 0.19   |
| Styrene                         | 0.005   | "     | <                                     | 0.20  | 88  | 0.22   |
| 1,1,1,2-Tetrachloroethane       | 0.010   | 15    | <                                     | 0.22  | 100   | 0.22   |
| 1,1,2,2-Tetrachloroethane       | 0.010   | п     | <                                     | 0.15  | 61  | 0.14   |
| Tetrachloroethene               | 0.020   | н     | <                                     | 0.15  | 97  | 0.14   |
| Toluene                         | 0.005   | 11    | <                                     | 0.24  | 94  | 0.23   |

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|   | Client ID:<br>Zenon ID:<br>Date Sampled: |         | BH3<br>13-15<br>055069 96<br>96/12/12 | BH3<br>13-15<br>055069 96<br>96/12/12 | BH3<br>13-15<br>055069 96<br>96/12/12 | BH3<br>13-15<br>055069 96<br>96/12/12 |
|---|--|---------|---------------------------------------|---------------------------------------|---------------------------------------|---------------------------------------|
| Component                                   | MDL                                      | Units   |                                       | MS                                    | % recoveries                          | MSD                                   |
| 1,1,1-Trichloroethane                       | 0.005                                    |         | <                                     | 0.25                                  | 100                                   | 0.25                                  |
| 1,1,2-Trichloroethane                       | 0.010                                    | "       | <                                     | 0.23                                  | 94                                    | 0.23                                  |
| Trichloroethene                             | 0.005                                    | u       | <                                     | 0.31                                  | 130                                   | 0.31                                  |
| Vinvl Acetate                               | 0.010                                    | п       | <                                     | 0.10                                  | 41                                    | 0.072                                 |
| Vinyl Chloride                              | 0.010                                    | "       | <                                     | 0.24                                  | 95                                    | 0.21                                  |
| Xylenes(Total)                              | 0.005                                    | н       | <                                     | 0.68                                  | 94                                    | 0.68                                  |
| Surrogate Recoveries                        |  | %       |                                       |                                       |                                       |                                       |
| d4-1,2-Dichloroethane                       |  |         | 99                                    | 99                                    | 99                                    | 95                                    |
| d8-Toluene                                  |  |         | 95                                    | 99                                    | 99                                    | 101                                   |
| Bromofluorobenzene                          |  |         | 92                                    | 108                                   | 108                                   | 100                                   |
| Semi Volatiles via SW846 Method             | 18270                                    |         |                                       | • Qr 5.                               |                                       |                                       |
| Phenol                                      | 0.11                                     | mg/kg   | <0.22                                 | _                                     | -                                     | -                                     |
| Bis(2-chloroethyl)ether                     | 0.18                                     | "       | < 0.36                                | _                                     | -                                     | -                                     |
| 2-Chlorophenol                              | 0.27                                     | ц       | < 0.54                                | -                                     | -                                     | _                                     |
| 1,3-Dichlorobenzene                         | 0.20                                     | н       | < 0.40                                | -                                     | -                                     | _                                     |
| 1-4-Dichlorobenzene                         | 0.20                                     | н       | < 0.40                                | -                                     | -                                     | _ 、                                   |
| 1,2-Dichlorobenzene                         | 0.20                                     | н       | <0.40                                 | _                                     | -                                     | -                                     |
| Bis(2-chloroisopropyl)ether                 | 0.15                                     | n       | < < 0.30                              | _                                     | -                                     | -                                     |
| Hexachloroethane                            | 0.20                                     | ц       | < 0.40                                | _                                     | _                                     | _                                     |
| N-Nitroso-di-N-Propylamine                  | 0.20                                     | 11      | <0.42                                 | _                                     | -                                     | -                                     |
| Nitrobenzene                                | 0.20                                     | "       | <0.42                                 | _                                     | -                                     | -                                     |
| lsophorone                                  | 0.20                                     | n       | < 0.80                                | -                                     | _                                     | -                                     |
| 2-Nitrophenol                               | 0:14                                     | н       | < 0.28                                | _                                     | -                                     |                                       |
| 2,4-Dimethylphenol                          | 0.17                                     |         | < 0.34                                | _                                     | _                                     | -                                     |
| Bis(2-chloroethoxy)methane                  | 0.13                                     | п       | <0.26                                 | _                                     | -                                     | _                                     |
| 2,4-Dichlorophenol                          | 0.12                                     | 11      | <0.20                                 | _                                     | -                                     | _                                     |
| 1,2.4-Trichlorobenzene                      | 0.12                                     | "       | <0.24                                 | -                                     | _                                     | _                                     |
| Naphthalene                                 | 0.03                                     |         | <0.40                                 | -                                     | _                                     | -                                     |
| Hexachlorobutadiene                         | 0.20                                     | и       | <0.40                                 | -                                     | -                                     | -                                     |
| 4-Chloro-3-Methylphenol                     | 0.14                                     | "       | <0.40                                 |                                       | _                                     | -                                     |
| Hexachlorocyclopentadiene                   | 0.14                                     | "       | <0.28                                 | -                                     |                                       | -                                     |
| 2,4,6-Trichlorophenol                       | 0.20                                     | 9       | <0.40                                 | -                                     |                                       | _                                     |
| 2,4,0-11 chorophenor<br>2-Chloronaphthalene | 0.12                                     | н       | <0.24                                 | -                                     | _                                     |                                       |
| -   | 0.09                                     | 11      | <0.18                                 | -                                     | -                                     | -                                     |
| Accenaphthylene                             | 0.04                                     | 11      | <0.08                                 | -                                     | -                                     | -                                     |
| Dimethyl phthalate<br>2,6-Dinitrotoluene    | 0.06                                     | 11      | <0.22                                 | -                                     | -                                     | _                                     |
|   | 0.08                                     | *1      | <0.12                                 | -                                     | _                                     | _                                     |
| Acenaphthene                                | 0.48                                     | u       | <0.14<br><0.96                        | -                                     | -                                     | -                                     |
| 2,4-Dinitrophenol<br>2,4-Dinitrotoluene     | 0.48                                     | "       | <0.96<br><0.10                        | -                                     | -                                     | -                                     |
| 4-Nitrophenol                               | 0.14                                     | "       | <0.10                                 | -                                     | -                                     | -                                     |
| 1   |  | 0       |                                       | -                                     | -                                     | -                                     |
| Fluorene                                    | 0.03                                     | н       | < 0.06                                | -                                     | -                                     | -                                     |
| 4-Chlorophenylphenylether                   | 0.09                                     |         | <0.18                                 | -                                     | -                                     | -                                     |
| Diethyl phthalate                           | . 0.11                                   | u       | <0.22                                 | -                                     | -                                     | -                                     |
| 4,6-Dinitro-2-methylphenol                  | 0.15                                     |         | < 0.30                                | -                                     | -                                     | -                                     |
| N-Nitrosodiphenylamine                      | 0.19                                     | "       | < 0.38                                | -                                     | -                                     | -                                     |
| 4-Bromophenylphenylether                    | 0.03                                     | e<br>11 | < 0.06                                | -                                     | -                                     | -                                     |
| Hexachlorobenzene                           | 0.20                                     | n       | <0.40                                 | -                                     | -                                     | -                                     |

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|                            | ~             |       | BH3       | BH3       | BH3          | BH3       |
|----------------------------|---------------|-------|-----------|-----------|--------------|-----------|
|                            | Client ID:    |       | 13-15     | 13-15     | 13-15        | 13-15     |
|                            | Zenon ID:     |       | 055069 96 | 055069 96 | 055069 96    | 055069 96 |
|                            | Date Sampled: |       | 96/12/12  | 96/12/12  | 96/12/12     | 96/12/12  |
| Component                  | MDL           | Units |           | MS        | % recoveries | MSD       |
| Pentachlorophenol          | 0.11          | "     | <0.22     | -         | -            | -         |
| Phenanthrene               | 0.03          | 11    | < 0.06    | -         | -            | -         |
| Anthracene                 | 0.02          | "     | < 0.04    | -         | -            | -         |
| Di-n-butyl phthalate       | 0.11          | +1    | <0.22     | -         | -            | -         |
| Fluoranthene               | 0.02          | 11    | <0.04     | -         | -            | -         |
| Pyrene                     | 0.03          | п     | <0.06     | -         | -            | -         |
| Benzyl butyl phthalate     | 0.06          | 11    | < 0.12    | -         | -            | -         |
| 3,3-Dichlorobenzidine      | 0.10          | н     | <0.20     | -         | -            | -         |
| Benzo(a)anthracene         | 0.02          | "     | < 0.04    | -         | -            | -         |
| Chrysene                   | 0.03          | "     | < 0.06    | -         | -            | -         |
| Bis(2-ethylhexyl)phthalate | 0.14          | н.,   | <0.28     | -         | -            | -         |
| Di-n-octyl phthalate       | 0.11          | "     | < 0.22    | -         | -            | -         |
| Benzo(b)fluoranthene       | 0.04          | "     | < 0.08    | -         | -            | -         |
| Benzo(k)fluoranthene       | 0.04          | 11    | < 0.08    | -         | -            | - 1       |
| Benzo(a)pyrene             | 0.05          | м     | <0.10     | -         | -            | -         |
| Indeno(1,2,3-cd)pyrene     | 0.06          | 11    | <0.12     | -         | -            |           |
| Dibenzo(a,h)anthracene     | 0.04          | "     | < 0.08    | -         | -            | -         |
| Benzo(ghi)perylene         | 0.04          | u.    | < 0.08    | -         | · _ ·        |           |
| N-Nitrosodimethylamine     | 1.0           | mg/kg | <2.0      | -         | -            | -         |
| Aniline                    | 0.50          | "     | <1.0      | _         | -            | -         |
| Benzyl alcohol             | 0.50          | "     | <1.0      | -         | -            | -         |
| Carbazole                  | 0.50          | "     | <1.0      | -         | -            | -         |
| 2-Methylphenol             | 0.50          | **    | <1.0      | -         | -            | -         |
| Benzoic acid               | 0.50          | "     | <1.0      | -         | -            | -         |
| 4-Chloroaniline            | 0.50          | ц     | <1.0      | -         | -            | -         |
| 2-Methylnaphthalene        | 0.10          | "     | < 0.20    | -         | -            | -         |
| 2,4,5-Trichlorophenol      | 0.10          | н     | < 0.20    | -         |              | -         |
| 2-Nitroaniline             | 0.50          | 51    | <1.0      | -         | -            | ~         |
| 3-Nitroaniline             | 0.50          | 74    | <1.0      | _         | -            | -         |
| Dibenzofuran               | 0.50          | ч     | <1.0      | -         | -            | -         |
| Benzidine                  | 0.50          | **    | <1.0      | -         |              | ~         |
| 4-Nitroaniline             | 0.50          | "     | <1.0      | -         | -            | -         |
| Surrogate Recoveries       |               | %     |           |           |              |           |
| 2-Fluorophenol             |               |       | 43        | -         | -            | -         |
| d5-Phenol                  |               |       | 68        | -         | -            | -         |
| d5-Nitrobenzene            |               |       | 46        | -         | _            | -         |
| 2-Fluorobiphenyl           |               |       | 65        | -         | -            | -         |
| 2,4,6-Tribromophenol       |               |       | 57        | _         | -            | -         |
| d14-p-Terphenyl            |               |       | 72        | _         | -            | -         |
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|                                 |               |          | BH3          |
|---------------------------------|---------------|----------|--------------|
|                                 | Client ID:    |          | 13-15        |
|                                 | Zenon ID:     |          | 055069 96    |
|                                 | Date Sampled: |          | 96/12/12     |
| Component                       | MDL           | Units    | % recoveries |
| Mercury via SW846 Method 7471   |               |          |              |
| Mercury                         | 0.04          | mg/kg    | -            |
| Metals via SW846 Method 6010    |               | 0.0      |              |
| Arsenic                         | 5.0           | mg/kg    | -            |
| Selenium                        | 10            | "        | -            |
| Barium                          | 0.1           | **       | -            |
| Cadmium                         | 0.2           | 11       | -            |
| Chromium                        | 5             | 11       | -            |
| Lead                            | 10            |          | -            |
| Silver                          | 0.5           | "        | -            |
| Volatiles via SW846 Method 8260 |               |          |              |
| Acetone                         | 0.030         | mg/kg    | 130          |
| Aerolein                        | 0.010         | "        | 73           |
| Aervlonitrile                   | 0.010         | н        | 78           |
| Benzene                         | 0.005         |          | 92           |
| Bromoform                       | 0.010         | н        | 100          |
| Bromomethane                    | 0.010         | 11       | 120          |
| 2-Butanone                      | 0.015         |          | 75           |
| Carbon Disulfide                | 0.010         | n        | 82           |
| Carbon Tetrachloride            | 0.010         | 11       | 100          |
| Chlorobenzene                   | 0.005         | ti.      | 92           |
| Chlorodibromomethane            | 0.005         |          | 100          |
| Chloroethane                    | 0.010         | *1       | 98           |
| 2-Chloroethylvinylether         | 0.010         | n        | 60           |
| Chlorotorm                      | 0.005         | 11       | 91 -         |
| Chloromethane                   | 0.010         | 71       | 94           |
| 1,2-Dichlorobenzene             | 0.005         | u        | 91           |
| 1,3-Dichlorobenzene             | 0.005         | "        | 86           |
| 1,4-Dichlorobenzene             | 0.005         | 11       | 88           |
| Dichlorobromomethane            | 0.005         | ц        | 100          |
| 1,1-Dichloroethane              | 0.005         | "        | 90           |
| 1,2-Dichloroethane              | 0.005         | п        | 91           |
| 1,1-Dichloroethene              | 0.010         | 11       | 89           |
| cis-1,2-Dichloroethene          | 0.010         | ,,       | 85           |
| trans-1,2-Dichloroethene        | 0.010         | н        | 85           |
| 1,2-Dichloropropane             | 0.005         |          | 97           |
| cis-1,3-Dichloropropene         | 0.005         | ч        | 92           |
| trans-1,3-Dichloropropene       | 0.005         | н        | 94           |
| Ethylbenzene                    | 0.005         |          | 90           |
| 2-Hexanone                      | 0.005         | n        | 90<br>70     |
| Methylene Chloride              | 0.010         | u        | 95           |
| 4-Methyl-2-Pentanone            |               |          |              |
| •                               | 0.010         | n        | 77           |
| Styrene                         | 0.005         | u        | 89<br>100    |
| 1,1,1,2-Tetrachloroethane       | 0.010         | <u>и</u> | 100          |
| 1,1,2,2-Tetrachloroethane       | 0.010         |          | 57           |
| Tetrachloroethene               | 0.020         | u        | 99<br>05     |
| Toluene                         | 0.005         |          | 95           |

|  |               |       | BH3          |
|--|---------------|-------|--------------|
|  | Client ID:    |       | 13-15        |
|  | Zenon ID:     |       | 055069 96    |
|  | Date Sampled: |       | 96/12/12     |
| Component                              | MDL           | Units | % recoveries |
| 1,1,1-Trichloroethane                  | 0.005         | "     | 100          |
| 1,1,2-Trichloroethane                  | 0.010         | н     | 93           |
| Trichloroethene                        | 0.005         | "     | 120          |
| Vinyl Acetate                          | 0.010         | н     | 28           |
| Vinyl Chloride                         | 0.010         | "     | 84           |
| Xylenes(Total)                         | 0.005         | 11    | 90           |
| Surrogate Recoveries                   |               | %     |              |
| d4-1,2-Dichloroethane                  |               |       | 95           |
| d8-Toluene                             |               |       | 101          |
| Bromofluorobenzene                     |               |       | 100          |
| Semi Volatiles via SW846 Method 82     | 270           |       |              |
| Phenol                                 | 0.11          | mg/kg | -            |
| Bis(2-chloroethyl)ether                | 0.18          | **    | -            |
| 2-Chlorophenol                         | 0.27          | 11    | -            |
| 1,3-Dichlorobenzene                    | 0.20          | **    | -            |
| 1-4-Dichlorobenzene                    | 0.20          | 11    | · -          |
| 1,2-Dichlorobenzene                    | 0.20          | u     | -            |
| Bis(2-chloroisopropyl)ether            | 0.15          |       | -            |
| Hexachloroethane                       | 0.20          | н     | -            |
| N-Nitroso-di-N-Propylamine             | 0.21          | "     | -            |
| Nitrobenzene                           | 0.20          | 0     | -            |
| Isophorone                             | 0.40          | "     | -            |
| 2-Nitrophenol                          | 0.14          | 11    | -            |
| 2,4-Dimethylphenol                     | 0.17          | ч     | -            |
| Bis(2-chloroethoxy)methane             | 0.13          | 11    | -            |
| 2,4-Dichlorophenol                     | 0.12          | **    | -            |
| 1,2,4-Trichlorobenzene                 | 0.20          | 11    | -            |
| Naphthalene                            | 0.03          | н     | -            |
| Hexachlorobutadiene                    | 0.20          | ч     | -            |
| 4-Chloro-3-Methylphenol                | 0.14          | "     | -            |
| Hexachlorocyclopentadiene              | 0.20          | ŧI    | -            |
| 2,4,6-Trichlorophenol                  | 0.12          |       | -            |
| 2-Chloronaphthalene                    | 0.09          | 11    | -            |
| Acenaphthylene                         | 0.04          | "     | -            |
| Dimethyl phthalate                     | 0.11          | ц     | -            |
| 2,6-Dinitrotoluene                     | 0.06          | 11    | -            |
| Acenaphthene                           | 0.07          | "     | -            |
| 2,4-Dinitrophenol                      | 0.48          | "     | -            |
| 2,4-Dinitrotoluene                     | 0.05          | u     | -            |
| 4-Nitrophenol                          | 0.14          | н     | -            |
| Fluorene                               | 0.03          | u     | -            |
| 4-Chlorophenylphenylether              | 0.09          | ц     | _            |
| Diethyl phthalate                      | 0.11          | и     | -            |
| 4,6-Dinitro-2-methylphenol             | 0.15          | н     | -            |
| N-Nitrosodiphenylamine                 | 0.19          | и     | -            |
| 4-Bromophenyiphenylether               | 0.03          | н     | -            |
| ······································ | 0.01          |       |              |

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|                            |               |       | BH3          |
|----------------------------|---------------|-------|--------------|
|                            | Client ID:    |       | 13-15        |
|                            | Zenon ID:     |       | 055069-96    |
|                            | Date Sampled: |       | 96/12/12     |
| Component                  | MDL           | Units | % recoveries |
| Pentachlorophenol          | 0.11          | "     | -            |
| Phenanthrene               | 0.03          | n     | -            |
| Anthracene                 | 0.02          | **    | -            |
| Di-n-butyl phthalate       | 0.11          | 0     | -            |
| Fluoranthene               | 0.02          | 9     | -            |
| Pyrene                     | 0.03          |       | -            |
| Benzyl butyl phthalate     | 0.06          | **    | -            |
| 3,3-Dichlorobenzidine      | 0.10          | a     | -            |
| Benzo(a)anthracene         | 0.02          | "     | -            |
| Chrysene                   | 0.03          | "     | -            |
| Bis(2-ethylhexyl)phthalate | 0.14          | ш     | -            |
| Di-n-octyl phthalate       | 0.11          |       | -            |
| Benzo(b)fluoranthene       | 0.04          | 11    | -            |
| Benzo(k)fluoranthene       | 0.04          | **    | -            |
| Benzo(a)pyrene             | 0.05          | 11    | -            |
| Indeno(1,2,3-cd)pyrene     | 0.06          | "     | -            |
| Dibenzo(a,h)anthracene     | 0.04          | 11    | -            |
| Benzo(ghi)perylene         | 0.04          | "     | -            |
| N-Nitrosodimethylamine     | 1.0           | mg/kg | -            |
| Aniline                    | 0.50          | "     | -            |
| Benzyl alcohol             | 0.50          | 11    | -            |
| Carbazole                  | 0.50          | н     | -            |
| 2-Methylphenol             | 0.50          | 11    | -            |
| Benzoie acid               | 0.50          | **    | -            |
| 4-Chloroaniline            | 0.50          | a     | -            |
| 2-Methylnaphthalene        | 0.10          | 0     | -            |
| 2,4,5-Trichlorophenol      | 0.10          | 15    | -            |
| 2-Nitroaniline             | 0.50          | **    | -            |
| 3-Nitroaniline             | 0.50          |       | -            |
| Dibenzofuran               | 0.50          | **    | -            |
| Benzidine                  | 0.50          | 1)    | -            |
| 4-Nitroaniline             | 0.50          | n     | -            |
| Surrogate Recoveries       |               | %     |              |
| 2-Fluorophenol             |               |       | -            |
| d5-Phenol                  |               |       | -            |
| d5-Nitrobenzene            |               |       | -            |
| 2-Fluorobiphenyl           |               |       | -            |
| 2,4,6-Tribromophenol       |               |       | -            |
| d14-p-Terphenyl            |               |       | -            |
|                            |               |       |              |

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## ZEL Summary of Analysis Pre. Dates

| Batch Code:                       | 1220ASA1        |
|-----------------------------------|-----------------|
| Mercury                           | 055066 96       |
|                                   | 055067 96       |
|                                   | 055068 96       |
|                                   | 055069 96       |
| Date analysed:                    | 96/12/20        |
| Date prepared:                    | 96/12/20        |
| Batch Code:                       | 1223ASC1        |
| Metals                            | 055066 96       |
|                                   | 055067 96       |
|                                   | 055068 96       |
|                                   | 055069 96       |
| Date analysed:                    | 96/12/30        |
| Date prepared:                    | 96/12/23        |
| Potek Code                        | 122058401       |
| Batch Code:<br>Volatiles via 8240 | 1220SM01        |
| volatiles via 8240                | 055066 96       |
|                                   | 055067 96       |
|                                   | 055068 96       |
| Data analyzadi                    | 055069 96       |
| Date analysed:                    | 96/12/20        |
| Date prepared:                    | 96/12/20        |
| Batch Code:                       | 1217PB02        |
| Volatiles via 8270                | 055066 96       |
|                                   | 055067 96       |
|                                   | 055068 96       |
|                                   | 055069 96       |
| Date analysed:                    | 96/12/20        |
| Date prepared:                    | 96/12/20        |
| F                                 | 2 \2) • +#/ 1 / |

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| COC Serial No. C 3017   | JSDAG Comments  | e Date Time<br>(1012) 17:30p-      | Airbill No. 748387340)<br>Fax & mail result ×0  |
|---|---|------------------------------------|---|
| n of Custody Record<br>oe Boad<br>0, NM 87401<br>(505) 326-2262 Phonend<br>(505) 326-2388 FAX | Type of Analysis and Both and | Received By:       Time       ISoD | dEX<br>Bonotes: Standard Turnaround<br>Chance @ Faunington  |
| Chain of C<br>termington, NM 87401  | Septic Leuch Fielb<br>Phase. Task 6006.77<br>Phase. Task 6006.77<br>Phase. Task 6006.77<br>Le lingthoy Antario and<br>Date Time Matrix<br>Date Time Matrix<br>Date Time Matrix<br>Date Time Matrix  | y:<br>Signature Date               | amples Iced:       Ves       No       Carrier:       F_e       E         servatives (ONLY for Water Samples)       Shipping and Lab Notes:       Cyanide       Shipping and Lab Notes:         Cyanide       Volatile Organic Analysis       Nitric acid (HNO3)       Shipping and Lab Notes:         Metals       Nitric acid (HNO3)       Nitric acid (HNO3)       Over Y       V         Other (Specify)       Other (Specify)       Over Y       CV       V |
| ) (197 - 1911)<br>197   | Project Name LFS<br>Project Number 17039<br>Samplers M Z<br>Laboratory Name Z<br>Sample Number (and depth)<br>BHJ-13-13 (06<br>BHJ-13-15 (06<br>BHJ-13-15 (06)  | Relinquished by:                   | Samples Iced: V Yes<br>Preservatives (ONLY for Water Samples)<br>Cyanide Solution Sodiu<br>Volatile Organic Analysis Solution Solution<br>Metals Specify Specify Solution Solution Solution Solution (Specify)  |

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5555 North Service Road Burlington, Ontario, Canada L7L 5H7 Tel: (905) 332-8788 Fax: (905) 332-9169

# Certificate of Analysis

#### CLIENT INFORMATION

### LABORATORY INFORMATION

| Attention:    | Cory Chance               | Contact:        | Ada Blythe, B.Sc., C.Chem. |
|---------------|---------------------------|-----------------|----------------------------|
| Client Name:  | Philip Environmental Inc. | Project:        | AN961368                   |
| Project:      | 17039                     | Date Received:  | 97/01/27                   |
| Project Desc: | WFS Septic Leach Field    | Date Reported:  | 97/02/04                   |
| Address:      | 4000 Monroe Road          | Submission No.: | 7A0483                     |
|               | Farmington, NM            | Sample No.:     | -002309                    |
|               | 87401                     |                 |                            |
| Fax Number:   | 505 326-2388              |                 |                            |
| Phone Number: | 505 326-2262              |                 |                            |

NOTES: '-' = not analysed '<' = less than Method Detection Limit (MDL) 'NA' = no data available LOQ can by determined for all analytes by multiplying the appropriate MDL X 3.33 Solids data is based on dry weight except for biota analyses. Organic analyses are not corrected for extraction recovery standards except for isotope dilution methods, (i.e. CARB 429 PAH, all PCDD/F and DBD/DBF analyses)

Methods used by Zenon are based upon those found in 'Standard Methods for the Examination of Water and Wastewater', Seventeenth Edition. Other methods are based on the principles of MISA or EPA methodologies.

All work recorded herein has been done in accordance with normal professional standards using accepted testing methodologies, quality assurance and quality control procedures except where otherwise agreed to by the client and testing company in writing. Any and all use of these test results shall be limited to the actual cost of the pertinent analysis done. There is no other warranty expressed or implied. Your samples will be retained at Zenon for a period of three weeks from receipt of data or as per contract.

COMMENTS:

"\*" Suspect chloride interference on silver recovery

"\*" There are no recoveries of 2-chlorovinylether, as this compound degrades very rapidly at room temperature. It can degrade in the time elapsing from sample set up to the time it is actually injected on the GC/MS.

Certified by:

Page 1

| Arsenic (gfaa) 0.0                                       | Client ID:<br>Zenon ID:<br>Date Sampled:<br>Component MDL (<br>Arsenic via SW846 Method 7060 | Units | Method<br>Blank<br>002309 97<br>97/01/24 | Blank<br>Spike<br>002309 97<br>97/01/24 | Blank<br>Spike<br>002309 97<br>97/01/24<br>%Recovery | WFS<br>ST-12497<br>002311-97<br>97/01/24 | WFS<br>ST-12497<br>002311 97<br>97/01/24<br>Duplicate | WFS<br>ST-12497<br>002311 97<br>97/01/24<br>M. Spike | WFS<br>ST-12497<br>002311 97<br>97/01/24<br>MS % Rec. | WFS<br>ST-12497<br>002311 97<br>97/01/24<br>MS Dup | WFS<br>ST-12497<br>(0(2311-97<br>97/01/24<br>MSD % Rec. |
|--|--|-------|--|---|--|--|---|--|---|--|---|
|  | _  | mg/L  | $\vee$                                   | 0.049                                   | 06   | V  | V   | ().()47  | 86  | 0.043  | 78  |
| Selenium via SW846 Method 7740<br>Selenium (gfaa) 0.0020 |  | mg/L  | V  | 0.050                                   | 89   | V  | 0.0026  | 0.045  | 78  | 0.043  | 74  |
| Metals via SW846 Method 6010                             | 010  |       |  |   |  |  |   |  |   |  |   |
| um   | _  | mg/L  | <0.033                                   | 2.3                                     | 110  | 0.26                                     | 0.33  | 2.7  | 011   | 2.6  | 100   |
|  | 0.001  | ÷     | 0.002                                    | 1.1                                     | 100  | 0.49                                     | 0.45  | 1.6  | 100   | 1.6  | 100   |
| Beryllium 0.0  | 0.001  | :     | V  | 0.58                                    | 110  | $\vee$                                   | V   | 0.61   | 110   | 0.59   | 110   |
| Cadmium 0.0  | 0.002  | ÷     | 0.004                                    | 0.56                                    | 100  | 0.004                                    | 0.004   | 0.59   | 110   | 0.57   | 100   |
| Calcium 0.   | 0.20   | :     | <0.22                                    | 11                                      | 100  | 39                                       | 38  | 48   | 91  | 47   | 83  |
| Chromium 0.0   | 0.004  | ÷     | $\vee$                                   | 1.1                                     | 100  | $\vee$                                   | V   | 1.2  | 110   | 1.1  | 100   |
| Cobalt 0.0   | 0.010  | 5     | <0.011                                   | 1.1                                     | 100  | <0.011                                   | <0.011  | 1.2  | 110   | 1.1  | 100   |
| Copper 0.0   | 0.005  | z     | <0.006                                   | 1.1                                     | 100  | 0.11                                     | 0.10  | 1.3  | 110   | 1.3  | 100   |
| Iron 0.0   | 0.010  | E     | <0.011                                   | 14                                      | 100  | 0.38                                     | 0.43  | 15   | 110   | 14   | 110   |
| Lead 0.0   | 020  | Ŧ     | 0.023                                    | 1.1                                     | 100  | <0.022                                   | <0.022  | 1.2  | 110   | 1.2  | 110   |
| Magnesium 0.0  | 0.050  | :     | <0.055                                   | 12                                      | 100  | 7.5                                      | 7.3   | 20   | 100   | 20   | 100   |
| Manganese 0.0  | 900  | ÷     | <0.007                                   | 1.1                                     | 100  | 0.031                                    | 0.035   | 1.2  | 110   | 1.2  | 100   |
| Molybdenum 0.0   | 0.010  | ÷     | <0.011                                   | 0.55                                    | 100  | 0.017                                    | <0.011  | 0.59   | 110   | 0.58   | 100   |
| Nickel 0.0   | 010  | Ŧ     | <0.011                                   | 0.57                                    | 100  | <0.011                                   | <0.011  | 0.59   | 110   | 0.59   | 110   |
| Phosphorus 0.0   | 0.060  | -     | <0.066                                   | 5.8                                     | 110  | 8.6                                      | 8.4   | 14   | 100   | 41   | 95  |
| Potassium 1.0  | 1.000  | 2     | 1.1                                      | 12                                      | 66   | 23                                       | 22  | 34   | 66  | 34   | 100   |
|  | 0.010  | ÷     | <0.011                                   | 0.57                                    | 100  | <0.011                                   | <0.011  | 0.29   | 53*   | 0.38   | 70*   |
| Sodium 0.  | 0.100  | :     | <0.11                                    | 11                                      | 100  | 44                                       | 43  | 54   | 96  | 53   | 87  |
| Thallium 0.0   | 0.060  | Ŧ     | <0.066                                   | 1.1                                     | 100  | <0.066                                   | <0.066  | 1.2  | 110   | 1.2  | 100   |
| Vanadium 0.0   | 0.005  | ÷     | <000.05                                  | 0.55                                    | 1()()  | <0.0010>                                 | <0.006  | 0.56   | 100   | 0.55   | 100   |
| Zinc 0.  | 0.005  | £     | <0.006                                   | 2.3                                     | 100  | 0.21                                     | 0.15  | 2.4  | 100   | 2.4  | 66  |

Page 2 of 11

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| 2/4/97                                   |                       |        | Zenc      | on Enviro | onmental  | Laboratori | Zenon Environmental Laboratories - Certificate of Analysis | tate of Ana | lysis      |            | Page 3 of 11 |
|--|-----------------------|--------|-----------|-----------|-----------|------------|--|-------------|------------|------------|--------------|
|  |                       |        | Method    | Blank     | Blank     | TCLP -WFS  | TCLP - WFS   | TCLP -WFS   | TCLP - WFS | TCLP - WFS | TCLP - WFS   |
| CI                                       | Client ID:            |        | Blank     | Spike     | Spike     | STS-12497  | STS-12497  | STS-12497   | STS-12497  | STS-12497  | STS-12497    |
| Zei                                      | Zenon ID:             |        | 002309 97 | 002309 97 | 002309 97 | 002313 97  | 00231397   | 002313 97   | 002313 97  | 002313 97  | 002313 97    |
| Date Sa                                  | Date Sampled:         |        | 97/01/24  | 97/01/24  | 97/01/24  |            |  |             |            |            |              |
| Component MDL<br>Mercury via Method 7470 | MDL Units<br>hod 7470 | Units  |           |           | %Recovery |            | Duplicate  | M. Spike    | MS % Rec.  | MS Dup     | MSD % Rec.   |
| Mercury(cvaa)                            | 0.05                  | ug/L   | V         | 1.0       | 100       | V          | V  | 1.0         | 100        | 1.0        | 100          |
| TCLP Metals via SW846 Method 6010        | SW846 .               | Method | 0109      |           |           |            |  |             |            |            |              |
|  |                       | mg/L   | <0.55     | <0.55     | 95        | <0.55      | <0.55  | 0.61        | 110        | 0.61       | 110          |
|  | 0.001                 | ÷      | V         | 1.1       | 100       | 1.2        | 1.2  | 2.2         | 110        | 2.2        | 110          |
| Cadmium                                  | 0.002                 | z      | V         | 0.56      | 100       | 0.002      | V  | 0.58        | 110        | 0.58       | 110          |
| Chromium                                 | 0.004                 | ŧ      | V         | 1.1       | 100       | V          | V  | 1.1         | 100        | 1.1        | 100          |
| Lead                                     | 0.020                 | ÷      | <0.022    | 1.1       | 100       | <0.022     | <0.022   | 1.2         | 110        | 1.2        | 110          |
| Selenium                                 | 0.060                 | ÷      | <0.066    | 0.58      | 100       | <0.066     | <0.066   | 0.61        | 110        | 0.62       | 120          |
| Silver                                   | 0.010                 | ÷      | <0.011    | 0.56      | 100       | <0.011     | <0.011   | 0.54        | 66         | 0.53       | -76          |
|  |                       |        |           |           |           |            |  |             |            |            |              |

Client: Philip Environmental Inc. Project: 17039

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|                           | Client ID:<br>Zenon ID:<br>Date Sampled: |       | Method<br>Blank<br>002309 97<br>97/01/24 | Blank<br>Spike<br>002309 97<br>97/01/24 | Blank<br>Spike<br>002309 97<br>97/01/24 | WFS<br>ST-12497<br>002311 97<br>97/01/24 | WFS<br>ST-12497<br>002311 97<br>97/01/24 | WFS<br>ST-12497<br>002311 97<br>97/01/24 |
|---------------------------|--|-------|--|---|---|--|--|--|
| Component                 | MDL                                      | Units |  |   | %Recovery                               |  | M. Spike                                 | MS % Rec.                                |
| Volatiles via SW846 Me    |  | а     | 0.1                                      | 22                                      | 15                                      | 1.4                                      | 22                                       | 4.4                                      |
| Acetone                   | 4.0                                      | ug/L  | 8.1                                      | 22                                      | 45                                      | 14                                       | 22                                       | 44                                       |
| Acrolein                  | 2.0                                      |       | <  | 90                                      | 90                                      | <  | 85                                       | 85                                       |
| Acrylonitrile             | 2.0                                      |       | <  | 90                                      | 90                                      | <  | 84<br>44                                 | 84<br>87                                 |
| Benzene<br>Bromoform      | 1.0<br>2.0                               |       | <  | 49<br>50                                | 98                                      | <  | 44<br>45                                 | 87<br>91                                 |
| Bromotorm                 | 2.0                                      |       | <  | 30<br>45                                | 101<br>90                               | <  | 43<br>44                                 | 87                                       |
| 2-Butanone                | 3.0                                      |       | <  | 45<br>66                                | 90<br>112                               | <     <                                  | 44<br>52                                 | 100                                      |
| Carbon Disulfide          | 2.0                                      |       | < <                                      | 60<br>60                                | 112                                     | <  | 52<br>60                                 | 120                                      |
| Carbon Tetrachloride      | 2.0                                      |       |  | 50                                      | 120                                     | <  | 00<br>45                                 | 90                                       |
| Chlorobenzene             | 1.0                                      |       | < <                                      | 50<br>52                                | 100                                     | <  | 47                                       | 93                                       |
| Chlorodibromomethane      | 1.0                                      |       | <  | 52<br>50                                | 99                                      | <  | 43                                       | - 86                                     |
| Chloroethane              | 2.0                                      |       | <  | 45                                      | 99                                      | < .                                      | 46                                       | 92                                       |
| 2-Chloroethylvinylether   | 2.0                                      |       | <  | 51                                      | 102                                     | <  | *  | *  |
| Chloroform                | 1.0                                      |       | <  | 44                                      | 88                                      | 4.8                                      | 48                                       | 96                                       |
| Chloromethane             | 2.0                                      |       | <  | 49                                      | 98                                      | <  | 48                                       | 96                                       |
| 1.2-Dichlorobenzene       | 1.0                                      |       | <  | 50                                      | 100                                     | <  | 47                                       | 93                                       |
| 1,3-Dichlorobenzene       | 1.0                                      |       | <  | 52                                      | 103                                     | <  | 45                                       | 90                                       |
| 1,4-Dichlorobenzene       | 1.0                                      |       | <  | 53                                      | 105                                     | 37                                       | -80                                      | 92                                       |
| Dichlorobromomethane      | 1.0                                      | ••    | <  | 48                                      | 95                                      | <  | 43                                       | 87                                       |
| 1.1-Dichloroethane        | 1.0                                      |       | <  | 46                                      | 91                                      | <  | 45                                       | 89                                       |
| 1,2-Dichloroethane        | 1.0                                      |       | <  | 47                                      | 93                                      | <  | 45                                       | 89                                       |
| 1,1-Dichloroethene        | 2.0                                      |       | <  | 47                                      | 95                                      | <  | 45                                       | 90                                       |
| cis-1,2-Dichloroethene    | 2.0                                      |       | <  | 46                                      | 92                                      | <  | 45                                       | 90                                       |
| trans-1,2-Dichloroethene  | 2.0                                      |       | <  | 46                                      | 93                                      | <  | 45                                       | 91                                       |
| 1.2-Dichloropropane       | 1.0                                      | •     | <  | 49                                      | 97                                      | <  | 44                                       | 87                                       |
| cis-1,3-Dichloropropene   | 1.0                                      |       | <  | 49                                      | 98                                      | <  | 43                                       | 85                                       |
| trans-1,3-Dichloropropen  | ie 1.0                                   | • •   | <  | 48                                      | 95                                      | <  | 43                                       | 85                                       |
| Ethylbenzene              | 1.0                                      | ••    | <  | 51                                      | 102                                     | <  | 47                                       | 94                                       |
| 2-Hexanone                | 2.0                                      |       | <  | 52                                      | 103                                     | <  | 42                                       | 83                                       |
| Methylene Chloride        | 2.0                                      |       | 3.5                                      | 39                                      | 79                                      | 39                                       | 75                                       | 84                                       |
| 4-Methyl-2-Pentanone      | 2.0                                      | **    | <  | 51                                      | 101                                     | <  | 46                                       | 91                                       |
| Styrene                   | 1.0                                      | **    | <  | 52                                      | 105                                     | <  | 46                                       | 93                                       |
| 1,1,1,2-Tetrachloroethand |  | **    | <  | 51                                      | 103                                     | <  | 46                                       | 93                                       |
| 1,1,2,2-Tetrachloroethand |  | ,.    | <  | 53                                      | 107                                     | <  | 45                                       | 90                                       |
| Tetrachloroethene         | 4.0                                      | ••    | <  | 52                                      | 104                                     | <  | 47                                       | 93                                       |
| Toluene                   | 1.0                                      | ••    | <  | 52                                      | 104                                     | <  | 45                                       | 91                                       |
| 1,1,1-Trichloroethane     | 1.0                                      |       | <  | 51                                      | 103                                     | <  | 46                                       | 92                                       |
| 1,1,2-Trichloroethane     | 2.0                                      |       | <  | 48                                      | 97                                      | <  | 45                                       | 91                                       |
| Trichloroethene           | 1.0                                      | **    | <  | 52                                      | 103                                     | <  | 44                                       | 88                                       |
| Vinyl Acetate             | 2.0                                      | .,    | <  | 52                                      | 103                                     | <  | 56                                       | 110                                      |
| Vinyl Chloride            | 2.0                                      | **    | <  | 46                                      | 92<br>102                               | <  | 46                                       | 91                                       |
| Xylenes(Total)            | 1.0                                      | **    | <  | 154                                     | 103                                     | <  | 140                                      | 95                                       |
| Surrogate Recoveries      |  | %     | 00                                       | 0.1                                     | 0.1                                     | 0.1                                      | 0.1                                      | 0.5                                      |
| d4-1,2-Dichloroethane     |  |       | 92                                       | 96<br>104                               | 96                                      | 96                                       | 96                                       | 96                                       |
| d8-Toluene                |  |       | 99                                       | 104                                     | 104                                     | 100                                      | 100                                      | 100                                      |
| Bromofluorobenzene        |  |       | 92                                       | 105                                     | 105                                     | 94                                       | 103                                      | 103                                      |
| pН                        |  |       | 6.00                                     | -                                       | -                                       | 2.00                                     | -  | -  |

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|                          |               | WFS       | WFS        |
|--------------------------|---------------|-----------|------------|
|                          | Client ID:    | ST-12497  | ST-12497   |
|                          | Zenon ID:     | 002311 97 | 002311 97  |
|                          | Date Sampled: | 97/01/24  | 97/01/24   |
| Component                | MDL           | MS Dup    | MSD % Rec. |
| Volatiles via SW846 Me   |               |           |            |
| Acetone                  | 4.0           | 25        | 50         |
| Acrolein                 | 2.0           | 85        | 85         |
| Acrylonitrile            | 2.0           | 84        | 85         |
| Benzene                  | 1.0           | 43        | 86         |
| Bromoform                | 2.0           | 43        | 86         |
| Bromomethane             | 2.0           | 43        | 86         |
| 2-Butanone               | 3.0           | 48        | 96         |
| Carbon Disulfide         | 2.0           | 55        | 110        |
| Carbon Tetrachloride     | 2.0           | 43        | 85         |
| Chlorobenzene            | 1.0           | 46        | 91         |
| Chlorodibromomethane     | 1.0           | 43        | 86         |
| Chloroethane             | 2.0           | 43        | 87         |
| 2-Chloroethylvinylether  | 2.0           | *         | *          |
| Chloroform               | 1.0           | 46        | 91         |
| Chloromethane            | 2.0           | 45        | 90         |
| 1,2-Dichlorobenzene      | 1.0           | 47        | 94         |
| 1.3-Dichlorobenzene      | 1.0           | 46        | 83         |
| 1,4-Dichlorobenzene      | 1.0           | 78        | 90         |
| Dichlorobromomethane     | 1.0           | 42        | 84         |
| 1.1-Dichloroethane       | 1.0           | 42        | 85         |
| 1,2-Dichloroethane       | 1.0           | 43        | 86         |
| 1.1-Dichloroethene       | 2.0           | 43        | 85         |
| cis-1.2-Dichloroethene   | 2.0           | 44        | 87         |
| trans-1,2-Dichloroethene | 2.0           | 43        | 86         |
| 1.2-Dichloropropane      | 1.0           | 43        | 86         |
| cis-1.3-Dichloropropene  | 1.0           | 42        | 83         |
| trans-1,3-Dichloroproper | ne 1.0        | 41        | 83         |
| Ethylbenzene             | 1.0           | 46        | 91         |
| 2-Hexanone               | 2.0           | 41        | 83         |
| Methylene Chloride       | 2.0           | 72        | 80         |
| 4-Methyl-2-Pentanone     | 2.0           | 44        | 87         |
| Styrene                  | 1.0           | 45        | 90         |
| 1,1,1,2-Tetrachloroethan | e 2.0         | 46        | 92         |
| 1.1.2.2-Tetrachloroethan |               | 45        | 90         |
| Tetrachloroethene        | 4.0           | 45        | 89         |
| Toluene                  | 1.0           | 45        | 89         |
| 1,1,1-Trichloroethane    | 1.0           | 43        | 86         |
| 1,1,2-Trichloroethane    | 2.0           | 43        | 85         |
| Trichloroethene          | 1.0           | 43        | 85         |
| Vinyl Acetate            | 2.0           | 55        | 110        |
| Vinyl Chloride           | 2.0           | 41        | 82         |
| Xylenes(Total)           | 1.0           | 140       | 91         |
| Surrogate Recoveries     | 1.0           | 1-1-1-2   | 1 1        |
| d4-1,2-Dichloroethane    |               | 97        | 97         |
| d8-Toluene               |               | 99<br>99  | 99         |
| Bromofluorobenzene       |               | 102       | 102        |
| CITATION POLIZONO        |               |           | 102        |
| рН                       |               | -         | -          |
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| Component             | Client ID:<br>Zenon ID:<br>Date Sampled:<br>MDL | Units | Method<br>Blank<br>002309 97<br>97/01/24 | TCLP -WFS<br>STS-12497<br>002313 97 | TCLP -WFS<br>STS-12497<br>002313 97<br>Duplicate |
|-----------------------|---|-------|--|-------------------------------------|--|
| TCLP Volatiles via SW |   | Onits |  |                                     | Duplicate  |
| Benzene               | 0.2   | ug/L  | <2.0                                     | <2.0                                | <2.0   |
| 2-Butanone            | 5.0   | "     | <50                                      | <50                                 | <50  |
| Carbon Tetrachloride  | 0.3   | **    | <3.0                                     | <3.0                                | <3.0   |
| Chlorobenzene         | 0.6   |       | <6.0                                     | <6.0                                | <6.0   |
| Chloroform            | 0.4   | "     | <4.0                                     | <4.0                                | <4.()  |
| 1,2-Dichloroethane    | 0.4   |       | <4.0                                     | <4.()                               | <4.()  |
| 1,1-Dichloroethene    | 0.4   | **    | <4.0                                     | <4.0                                | <4.0   |
| Pyridine              | 250   | "     | <2500                                    | <2500                               | <2500  |
| Tetrachloroethene     | 0.2   | "     | <2.0                                     | <2.0                                | <2.0   |
| Trichloroethene       | 0.3   | 11    | <3.0                                     | <3.0                                | <3.0   |
| Vinyl Chloride        | 2.9   |       | <29                                      | <29                                 | <29  |
| 1,4-Dichlorobenzene   | 0.2   | 11    | <2.0                                     | 11                                  | 13   |
| Surrogate Recoveries  |   | %     |  |                                     |  |
| d4-1,2-Dichloroethane |   |       | 96                                       | 98                                  | 93   |
| d8-Toluene            |   |       | 101                                      | 95                                  | 95   |
| 1,4-Bromofluorobenzen | ie  |       | 101                                      | 100                                 | 96   |

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### Zenon Environmental Laboratories - Certificate of Analysis

Page 7 of 11

|  | Client ID:<br>Zenon ID:<br>ite Sampled: |       | Method<br>Blank<br>002309 97<br>97/01/24 | Blank<br>Spike 1<br>002309 97<br>97/01/24 | Blank<br>Spike 1<br>002309 97<br>97/01/24 | Blank<br>Spike 2<br>002309 97<br>97/01/24 | Blank<br>Spike 2<br>002309 97<br>97/01/24 | WFS<br>ST-12497<br>002311 97<br>97/01/24 |
|--|---|-------|--|---|---|---|---|--|
| Component  | MDL                                     | Units |  |   | %Recovery                                 |   | %Recovery                                 |  |
| Semi-Volatiles via SW846 M   |   | л     |  | 1.4                                       | 20  | 16  | 20  | 1.1                                      |
| Phenol<br>Di (2011 - di Di di  | 1.1                                     | ug/L  | <  | 14  | 28  | 16<br>NG                                  | 32  | 11                                       |
| Bis(2-chloroethyl)ether  | 1.8                                     |       | <  | NS<br>20                                  | -   | NS  | -   | <  |
| 2-Chlorophenol   | 2.7                                     | +1    | <  | 30<br>NS                                  | 61  | 36<br>NG                                  | 72  | <  |
| 1.3-Dichlorobenzene  | 2.0                                     | 11    | <  | NS  | -   | NS  | -   | <<br>20                                  |
| 1-4-Dichlorobenzene  | 2.0                                     |       | <  | 12  | 47  | 16<br>NG                                  | 64  | 20                                       |
| 1.2-Dichlorobenzene  | 2.0                                     | **    | <  | NS  | -   | NS  | -   | <  |
| Bis(2-chloroisopropyl)ether<br>Hexachloroethane  | 1.5<br>2.0`                             |       | <  | NS  | -   | NS<br>NS                                  | -   | <  |
|  |   | ,,    | <  | NS  | - 65                                      | 21  | 83  | <  |
| N-Nitroso-di-N-Propylamine<br>Nitrobenzene   | 2.1                                     |       | <  | 16<br>NS                                  | - 00                                      | NS  | 0.)<br>-                                  | <  |
| Isophorone   | 2.0<br>4.0                              |       | <  | NS  | -   | NS  | -   | <  |
| 2-Nitrophenol  | 4.0<br>1.4                              | *1    |  | NS  | -   | NS  | -   | <  |
| 2.4-Dimethylphenol   | 1.4                                     |       | <     <                                  | NS  | -   | NS  | -   | <  |
| Bis(2-chloroethoxy)methane   | - 1.3                                   |       | <  | NS  | -   | NS  | -   | <  |
| 2,4-Dichlorophenol   | 1.3                                     | ••    | <  | NS  | -   | NS  | -   | <  |
| 1,2,4-Trichlorobenzene   | 2.0                                     |       | <  | 12  | 49  | 17  | 67  | <  |
| Naphthalene  | 0.3                                     | ••    | <  | NS  | -   | NS  | -   | <  |
| Hexachlorobutadiene  | 2,0                                     | 14    | <  | NS  | -   | NS  | -   | <  |
| 4-Chloro-3-Methylphenol  | 2.0                                     |       | <  | 41  | 82  | 40  | 81  | <  |
| Hexachlorocyclopentadiene  | 2.0                                     | ••    | <  | 41<br>NS                                  | -   | 40<br>NS                                  | -   | <  |
| 2,4,6-Trichlorophenol  | 1,2                                     | .,    | <  | NS  | -   | NS  | -   | <  |
| 2Chloronaphthalene   | 0,9                                     |       | <  | NS  | -   | NS  | -   | <  |
| Acenaphthylene   | 0.4                                     | ,,    | <  | NS  | -   | NS  | -   | <  |
| Dimethyl phthalate   | 1.1                                     | 41    | <  | NS  | -   | NS  | _   | <  |
| 2,6-Dinitrotoluene   | 0.6                                     |       | <  | NS  | -   | NS  | _   | <  |
| Acenaphthene   | 0.0                                     |       | <  | 19  | -<br>76                                   | 21  | 83  | <  |
| 2,4-Dinitrophenol  | 4.8                                     | +1    | <  | NS  | -   | NS  | -   | <  |
| 2,4-Dinitrotoluene   | 4.8<br>0.5                              |       | <  | 21  | - 84                                      | 23  | 90  | <  |
| 4-Nitrophenol  | 1.4                                     |       | <  | 18  | 36  | 18  | 36  | <  |
| Fluorene   | 0.3                                     |       | <  | NS  |   | NS  | -   | <  |
| 4-Chlorophenylphenylether  | 0.9                                     |       | <  | NS  | -   | NS  | -   | <  |
| Diethyl phthalate  | 1.1                                     |       | <  | NS  | _   | NS  | -   | <  |
| 4,6-Dinitro-2-methylphenol   | 1.5                                     |       | <  | NS  | _   | NS  | _   | <  |
| N-Nitrosodiphenylamine   | 1.9                                     | н     | <  | NS  | -   | NS  | _   | <  |
| 4-Bromophenylphenylether   | 0.3                                     |       | <  | NS  | -   | NS  | _   | <  |
| Hexachlorobenzene  | 2.0                                     | ••    | <  | NS  | _   | NS  | _   | <  |
| Pentachlorophenol  | 1.1                                     |       | <  | 52  | 100                                       | 52  | 100                                       | <  |
| Phenanthrene   | 0.3                                     |       | <  | NS  | -   | NS  | -   | <  |
| Anthracene   | 0.2                                     | ••    | ~  | NS  | -   | NS  | _   | <  |
| Di-n-butyl phthalate   | 1.1                                     | ,,    | 1.4                                      | NS  | -   | NS  | -   | <  |
| Fluoranthene   | 0.2                                     |       | 1.4<br><                                 | NS  | -   | NS  | -   | <  |
| Pyrene   | 0.2                                     |       | <  | 24  | 96  | 24  | 95  | <  |
| Benzyl butyl phthalate   | 0.5                                     | ,,    | <  | 24<br>NS                                  | - 90                                      | 24<br>NS                                  | -   | 1.0                                      |
| 3,3-Dichlorobenzidine  | 1.0                                     | н     | <  | NS  | -   | NS  | -   | <  |
| Benzo(a)anthracene   | 0.2                                     |       | <  | NS  | -   | NS  | _   | <  |
| Chrysene   | 0.2                                     |       | <  | NS<br>NS                                  | -   | NS<br>NS                                  | -   | <  |
| Bis(2-ethylhexyl)phthalate   | 0.5<br>1.4                              |       | <  | NS<br>NS                                  | -   | NS<br>NS                                  | -   | 2.1                                      |
| Di-n-octyl phthalate   |   | .,    |  | NS<br>NS                                  |   | NS<br>NS                                  | -   | <  |
| Benzo(b)fluoranthene   | 1.1<br>0.4                              | .,    | < <                                      | NS<br>NS                                  | -   | INS<br>NS                                 | -   | <  |
| consector municipalities of the sector of th | U. <del>'</del> †                       |       |  | INO                                       | -   | UND                                       | -   |  |

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# Zenon Environmental Laboratories - Certificate of Analysis

Page 8 of 11

|                        | Client ID:    |       | Method    | Blank     | Blank     | Blank     | Blank     | WFS       |
|------------------------|---------------|-------|-----------|-----------|-----------|-----------|-----------|-----------|
|                        | Zenon ID:     |       | Blank     | Spike 1   | Spike 1   | Spike 2   | Spike 2   | ST-12497  |
|                        |               |       | 002309 97 | 002309 97 | 002309 97 | 002309 97 | 002309 97 | 002311.97 |
|                        | Date Sampled: |       | 97/01/24  | 97/01/24  | 97/01/24  | 97/01/24  | 97/01/24  | 97/01/24  |
| Component              |               | Units |           |           | %Recovery |           | %Recovery |           |
| Benzo(k)fluoranthene   | 0.4           | "     | <         | NS        | -         | NS        | -         | <         |
| Benzo(a)pyrene         | 0.5           |       | <         | NS        | -         | NS        | ~         | <         |
| Indeno(1,2,3-cd)pyrene | 0.6           | .,    | <         | NS        | -         | NS        | -         | <         |
| Dibenzo(a,h)anthracene | 0,4           | "     | <         | NS        | -         | NS        | -         | <         |
| Benzo(ghi)perylene     | 0.4           |       | <         | NS        | -         | NS        | -         | <         |
| N-Nitrosodimethylamine | 10            | **    | <         | NS        | -         | NS        | -         | <         |
| Aniline                | 5.0           |       | <         | NS        | -         | NS        | -         | <         |
| Carbazole              | 5.0           | "     | <         | NS        | -         | NS        | -         | <         |
| Benzyl alcohol         | 2.0           | **    | <         | NS        | -         | NS        | -         | <         |
| 2-Methylphenol         | 1.7           | н     | <         | NS        | -         | NS        | ~         | <         |
| Benzoic acid           | 1.0           | "     | <         | NS        | -         | NS        | -         | 620       |
| 4-Chloroaniline        | 5.0           |       | <         | NS        | -         | NS        | -         | <         |
| 2-Methylnaphthalene    | 1.0           | "     | <         | NS        | -         | NS        | -         | <         |
| 2,4,5-Trichlorophenol  | 0.6           | **    | <         | NS        | -         | NS        | -         | <         |
| 2-Nitroaniline         | 5.0           | 0     | <         | NS        |           | NS        | -         | <         |
| 3-Nitroaniline         | 5.0           |       | <         | NS        | -         | NS        | -         | <         |
| Dibenzofuran           | 5.0           | + 1   | <         | NS        | -         | NS        | -         | <         |
| Benzidine              | 10            | 0     | <         | NS        | -         | NS        | -         | <         |
| 4-Nitroaniline         | 5.0           | ••    | <         | NS        | -         | NS        | -         | <         |
| Surrogate Recoveries   |               | %     |           |           |           |           |           |           |
| 2-Fluorophenol         |               |       | 91        | 39        | 39        | 40        | 40        | 47        |
| d5-Phenol              |               |       | 64        | 29        | 29        | 31        | 31        | 35        |
| d5-Nitrobenzene        |               |       | 82        | 69        | 69        | 84        | 84        | 82        |
| 2-Fluorobiphenvl       |               |       | 77        | 70        | 70        | 79        | 79        | 73        |
| 2,4.6-Tribromophenol   |               |       | 95        | 98        | 98        | 92        | 92        | 83        |
| d-14-p-Terphenyl       |               |       | 93        | 98        | 98        | 95        | 95        | 76        |

Page 9 of 11

Zenon Environmental Laboratories - Certificate of Analysis

| Method         Blank         Blank         Spike 1           Client ID:         D02309 97         002309 97         002309 97         0           Date Sampled:         97/01/24         97/01/24         4 | Client ID:<br>Zenon ID:<br>Date Sampled: |              | Method<br>Blank<br>002309 97<br>97/01/24 | Blank<br>Spike 1<br>002309 97<br>97/01/24 | Blank<br>Spike 1<br>002309 97<br>97/01/24 | Blank<br>Spike 2<br>002309 97<br>97/01/24 | Blank<br>Spike 2<br>002309 97 | TCLP -WFS<br>STS-12497<br>002313 97 |
|---|--|--------------|--|---|---|---|-------------------------------|-------------------------------------|
| Component<br>TCLP Semi-Volatiles vi   | MDL<br>ia SW846 Method 8                 | Units<br>270 |  |   | %Recovery                                 |   | 7,Recovery                    |                                     |
| o-Cresol  | 1.7                                      | ug/L         | V  | NS  | ı   | NS  | ı                             | V                                   |
| m&p-Cresol  | 3.5                                      | :            | V  | 28  | 57  | 28  | 56                            | V                                   |
| 1,4-Dichlorobenzene   | 2.0                                      | -            | V  | 36  |   | 34  | 67                            | 7.4                                 |
| 2,4-Dinitrotoluene  | 0.5                                      | Ŧ            | V  | 48  |   | 51  | 100                           | V                                   |
| Nitrobenzene  | 2.0                                      | z            | V  | 43  |   | 41  | 83                            | V                                   |
| Pentachlorophenol   | 1.1                                      | ÷            | V  | 61  |   | 66  | 130                           | V                                   |
| 2.4.5-Trichlorophenol   | 0.6                                      | :            | V  | NS  |   | NS  | ł                             | V                                   |
| 2,4,6-Trichlorophenol   | 1.2                                      | ŧ            | V  | 44  |   | 45  | 06                            | V                                   |
| Hexachloroethane  | 2.0                                      | =            | V  | 31  |   | 28  | 57                            | V                                   |
| Hexachlorobutadiene   | 2.0                                      | ÷            | V  | 38  |   | 36  | 72                            | V                                   |
| Hexachlorobenzene   | 2.0                                      | ÷            | V  | 48  |   | 49  | 66                            | V                                   |
| Surrogate Recoveries  |  | $\gamma_c$   |  |   |   |   |                               |                                     |
| d5-Phenol   |  |              | 38                                       | 35  |   | 34  | 34                            | 52                                  |
| d5-Nitrobenzene   |  |              | 06                                       | 81  |   | 76  | 76                            | 91                                  |
| 2-Fluorobiphenyl  |  |              | 87                                       | 85  | 85  | 84  | 84                            | 92                                  |
| 2,4.6-Tribromophenol  |  |              | 94                                       | 92  | 92  | 79  | 67                            | 96                                  |
| d-14-p-Terphenyl  |  |              | 96                                       | 68  | 89  | 96  | 96                            | 79                                  |

Client: Philip Environmental Inc. Project: 17039

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| Batch Code:     | 0130ASB1  |           |
|-----------------|-----------|-----------|
| Arsenic (gfaa)  | 002309 97 |           |
|                 | 002311 97 |           |
| Date analysed   | 97/01/30  |           |
| Date prepared   | 97/01/30  |           |
| Batch Code:     | 0128ASA2  | 0130ASC1  |
| Mercury         | 002309 97 | 002313 97 |
|                 | 002311 97 |           |
| Date analysed   | 97/01/28  | 97/01/30  |
| Date prepared   | 97/01/28  | 97/01/30  |
| Batch Code:     | 0130ASB1  |           |
| Selenium (gfaa) | 002309 97 |           |
|                 | 002311 97 |           |
| Date analysed   | 97/01/30  |           |
| Date prepared   | 97/01/30  |           |
| Batch Code:     | 0130ASB1  |           |
| ICP Metals      | 002309 97 |           |
|                 | 002311 97 |           |
| Date analysed   | 97/01/30  |           |
| Date prepared   | 97/01/30  |           |
| Batch Code:     | 0130ASD1  |           |
| TCLP Metals     | 002309 97 |           |
|                 | 002313 97 |           |
| Date analysed   | 97/01/30  |           |
| Date prepared   | 97/01/30  |           |
| Batch Code:     | 0128SM01  |           |
| Volatiles       | 002309 97 |           |
|                 | 002311 97 |           |
| Date analysed   | 97/01/28  |           |
| Date prepared   | 97/01/28  |           |
| Batch Code:     | 0128SM01  |           |
| pН              | 002309 97 |           |
|                 | 002311 97 |           |
| Date analysed   | 97/01/28  |           |
| Date prepared   | 97/01/28  |           |
| Batch Code:     | 0130SM01  |           |
| TCLP Volatiles  | 002309 97 |           |
|                 | 002313 97 |           |
| Date analysed   | 97/01/30  |           |
| Date prepared   | 97/01/30  |           |
|                 |           |           |

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Date prepared

| Batch Code:         | 0129JM01  |
|---------------------|-----------|
| Semi-Volatiles      | 002309 97 |
|                     | 002311 97 |
| Date analysed       | 97/01/31  |
| Date prepared       | 97/01/29  |
| Batch Code:         | 0129JM02  |
| TCLP Semi-Volatiles | 002309 97 |
|                     | 002313 97 |
| Date analysed       | 97/01/31  |

97/01/29

Client: Philip Environmental Inc. Project: 17039

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5555 North Service Road Burlington, Ontario, Canada L7L 5H7 Tel: (905) 332-8788 Fax: (905) 332-9169

LABORATORY INFORMATION

# Certificate of Analysis

### CLIENT INFORMATION

| Attention:    | Cory Chance               | Contact:        | Ada Blythe, B.Sc., C.Chem. |
|---------------|---------------------------|-----------------|----------------------------|
| Client Name:  | Philip Environmental Inc. | Project:        | AN961368                   |
| Project:      | 17039                     | Date Received:  | 97/01/27                   |
| Project Desc: | WFS Septic Leach Field    | Date Reported:  | 97/02/04                   |
|               |                           |                 |                            |
| Address:      | 4000 Monroe Road          | Submission No.: | 7A0483                     |
|               | Farmington, NM            | Sample No.:     | 002310-002312              |
|               | 87401                     |                 |                            |
| Fax Number:   | 505 326-2388              |                 |                            |
| Phone Number: | 505 326-2262              |                 |                            |

NOTES: ''-' = not analysed '<' = less than Method Detection Limit (MDL) 'NA' = no data available LOQ can by determined for all analytes by multiplying the appropriate MDL X 3.33 Solids data is based on dry weight except for biota analyses. Organic analyses are not corrected for extraction recovery standards except for isotope dilution methods, (i.e. CARB 429 PAH, all PCDD/F and DBD/DBF analyses)

Methods used by Zenon are based upon those found in 'Standard Methods for the Examination of Water and Wastewater', Seventeenth Edition. Other methods are based on the principles of MISA or EPA methodologies.

All work recorded herein has been done in accordance with normal professional standards using accepted testing methodologies, quality assurance and quality control procedures except where otherwise agreed to by the client and testing company in writing. Any and all use of these test results shall be limited to the actual cost of the pertinent analysis done. There is no other warranty expressed or implied. Your samples will be retained at Zenon for a period of three weeks from receipt of data or as per contract.

COMMENTS: Revised report

Certified by

Page 1

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| Component<br>TCLP Leachate Procedure  | Client ID:<br>Zenon ID:<br>Date Sampled:<br>MDL | Units  | Method<br>Blank<br>002310 97<br>97/01/24 | Blank<br>Spike<br>002310 97<br>97/01/24<br>M. Spike | Blank<br>Spike<br>002310 97<br>97/01/24<br>MS % Rcc. | WFS<br>STS-12497<br>002312 97<br>97/01/24 | WFS<br>STS-12497<br>002312 97<br>97/01/24<br>Duplicate | WFS<br>STS-12497<br>00231297<br>97/01/24<br>M. Spike | WFS<br>STS-12497<br>002312 97<br>97/01/24<br>MS % Rcc. |
|---|---|--------|--|---|--|---|--|--|--|
| pH alter 5.5 ml of 1N HCl addition<br>pH initial $(5g + 96.5m]$ water)<br>pH of extraction fluid (semi-vols/metals)<br>pH of extraction fluid (volatiles) |   |        |  |   |  | 1.80<br>8.25<br>4.95<br>4.95              |  | 1 1 1 1  |  |
| Soil pH r <i>ia SW846 9045</i>  |   |        | ·  | ŗ   | ·  | 7.85                                      | 7.79   | t  | Ţ  |
| Reactivity via ASTMD5058C<br>Reaction/flume/dust/gas<br>Temperature change (10 minutes)<br>Temperature change (initial)                                   |   | deg C  | , , ,                                    | 1 1 1   |  | 000                                       |  |  |  |
| Moisture  | 0.01  | (-)%)  | ,  | ·   | ı  | 31  | 1  | I  | ,  |
| Cyanide, total via Method 9010  | 0.100   | mg/kg  | V  | 0.49  | 98   | 0.14                                      | 1  | ľ  |  |
| Ignitability  | -   | mm/min | I  | I   | ſ  | V   | I  | ı  | ı  |
| Reactive Sulphide   | 0.2   | mg/kg  | $\vee$                                   | 8.4   | 87   | 260                                       | 290  | 1  | ŕ  |

Page 2 of 3

Zenon Environmental Laboratories - Certificate of Analysis

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2/10/97

Client: Philip Environmental Inc. Project: 17039

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| Batch Code:       | 0128SPA1  |
|-------------------|-----------|
| TCLP pH Readings  | 002312 97 |
| Date analysed     | 97/01/29  |
| Date prepared     | 97/01/28  |
| Batch Code:       | 0130MSA6  |
| Soil pH           | 002312 97 |
| Date analysed     | 97/01/30  |
| Date prepared     | 97/01/30  |
| Batch Code:       | 0129MSA1  |
| Reactivity        | 002312 97 |
| Date analysed     | 97/01/29  |
| Date prepared     | 97/01/29  |
| Batch Code:       | 0128SPA1  |
| Moisture          | 002312 97 |
| Date analysed     | 97/01/28  |
| Date prepared     | 97/01/28  |
| Batch Code:       | 0130MNA1  |
| Reactive Sulphide | 002310 97 |
|                   | 002312 97 |
| Date analysed     | 97/02/07  |
| Date prepared     | 97/02/07  |
| Batch Code:       | 0130R.JA1 |
| Cyanide total     | 002310 97 |
|                   | 002312 97 |
| Date analysed     | 97/01/30  |
| Date prepared     | 97/01/30  |
| Batch Code:       | 0128SPA1  |
| lgnitability      | 002312 97 |
| Date analysed     | 97/01/28  |
| 5                 | 97/01/28  |

N. L. F. L. المناعس s,∢)> いい & Cyanide analylis \*Selenium-GFAA, XXALL SUIFILE ďβ to reactivity herbicides on Time ر: د. Comments \*\*\* Do not Tax results to terricidus d Airbill No. 748387726 H9-CVAA 3020 T/LP  $\bigcirc$ Date 717 COC Serial No. 40(C>7 2 herit 2-40 AL LAT-X 1201 Shipping and Lab Notes: RUSH TURNAROUND. Ory Chance @ Fax # abour. E. 111 1501 10 T 15C 22 1VC/14 Signature \* (CHI 9011651 Received By: (505) 326-2262 Phone (505) 326-2388 FAX 10/20/ **Chain of Custody Record** Ì \*0109 A M 4-3 J ibxaly Type of Analysis and Bottle Time 1500 ЛX 4000 Monroe Road Farmington, NM 87401 Total Number of Bottles 6 Carrier: Hed Sludge Water 24/4×1 Water S/vdge Water Matrix Date Chance / K. Padilla 4751730 Phase . Task 600 .... Nitric acid (HNO3) .... Hydrochloric acid (HCI) 4~0 DDD ... Sodium hyroxide (NaOH) .. Sulfuric acid (H2SO4) Time 0001 0001 2101 Sigl +traw 76/42/ Å hario, Date -enon reservatives (ONLY for Water Samples) Chem > Yes ignature Sample Number (and depth) Land TPY LIZS-STS-12497 NES-575-12497 WFJ-JT-12497 WFJ-JJ-J2497 WFS-57-12497 Location ( 📈 Vorfatile Organic Analysis WFS Name elinquished by samplers  $\subset \mathcal{M}$ Project Number + Samples Iced: Other (Specify) 🔲 ТРН (418.1) ... WMetals ..... 1165 - 2210 િહ્યુદ્ધ - પ્રેર્ેો 🗌 🖒 Cyanide ... Project Name Laboratory Z

**APPENDIX C - CHEMICAL WASTE PROFILE** 

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| New D Amendment   |   | MATERIAL PROFILE NO<br>PXWFM-0001  |
|---|---|--|
| City/County_Bloon Field San J.<br>City/County_Bloon Field San J.<br>State_Ney_Mexice ZipCode<br>USEPAID#_ <u>PEHDIAC</u> _NM_98<br>State ID#<br>B. DOT Shipping Name <u>MOA</u> <u>PC</u><br><br>L. J. Q. J. D.   | 874D       City       FAAMIAGTR         16682-120       Attention       OR 9       C.         Telephone I       505       32.1         RA       UASTE       D. ANNUAL REPOR         fazard Class       SIC Code:       1         RQ       Source Code:       A         No       Process Generating:       Origin Code:       B         System Type:       M       M | 668     EXT.       JVIR.     (SOR761)       JWROE ROAD       A     State ALMZip Code 87401       B     Code 87401       B <t< th=""></t<>  |
| F. PHYSICAL CHARACTERISTICS AT 70° F         1: Infectious or Biological Waste? D Yes B No         2: NRC Regulated Badioactive? D Yes B No         3: Reactivity None       Water Reactive         D Pyrophoric       Shock Sensitive         D Cyanides       D OT Explosive         D Sulfides       D Other         I Gas (Cylinder)       Solid       %         Lab-Pack       Free Liquids       100 %         Layers       Single Layered       Bi-layered       Multi-layered         Viscosity       I Medium       Hich       Odor         None       Mild       Strong       Describe: | H. PHYSICALCHEMICAL CONSTITUENTS  | Dermal Toxicity LOs(Mg/Kg) $\Box \leq 40$ $\Box > 200, \leq 1000$ $\Box > 46, \leq 200$ $\boxtimes > 1000$ 4. Material poisonous by inhalation? $\Box$ YesOral Toxicity LDs(Mg/Kg) $\Box \leq 50$ $\Box > 5, \leq 50$ Solids: $\Box > 50, \leq 200$ Liquids: $\Box > 50, \leq 500$ Solids: $\Box > 50, \leq 500$ |
| G. METALS<br>ST NONE DTCLP (MG/L) DTOTAL (PPM)<br>Reg. Limit Betow Above Range<br>Arsenic 5 mg/L 0 0<br>Barium 100 mg/L 0 0<br>Cadmium 1 mg/L 0 0<br>Cadmium 5 mg/L 0 0<br>Copper 0 0<br>Lead 5 mg/L 0 0<br>Mercury 0.2 mg/L 0 0<br>Selenium 1 mg/L 0 0<br>Silver 5 mg/L 0 0<br>Cithers:  |   | I. ANTICIP TED VOLUME         Oty.       Centainer         Oty.       Centainer         Oty.       Centainer         Oty.       Sol, pall         Oty.       Super Sack*         Oty.       Super Sack*         Oty.       Sol, drum         Oty.       Rollot//Dump Trailer         Oty.       Sol, drum         Other       Sol, drum         Other       Other         Per       I Time         Vear       Other As point         Other As point       (1) Is this waste regulated as a Marine Pollutant  |

Generator's Certification: I hereby certify that the above and attached description is complete and accurate to the best of my knowledge and ability to determine that no deliberate or willful omissions of composition properties exist and that all known or suspected hazards have been disclosed. I certify that the materials tested are representative of all material described by this profile.

Generator's Authorized Signature:

Date 2/18/97

MILAGRO PLANT CHEMICAL LIST

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| ACID REAGENT            | 0.06  | Sulfamic Acid. Sodium Chloride   |
|-------------------------|-------|--|
| AMINO ACID              | 0.06  |  |
| AMINO ACID F            | 0.03  |  |
| CITRIC ACID             | 0.12  |  |
| FERROVER IRON           | 0.09  | Sodium Hydrosulfite, Sodium Bisulfite, Sodium Sulfite                    |
| HARDNESS INDICATOR      | 0.01  |  |
| HARDNESS BUFFER         | 0.01  | Borate, Tetra Sodium, penta hydrate, Potassium Hydroxide, Sodium Sulfide |
| .1 N HCL                | 0.49  |  |
| METHANOL                | 11.82 |  |
| MOLYBDATE POWDER        | 0.06  | Sodium Molybdate   |
| MOYLBDATE LIQUID        | 0.78  | Sulfuric Acid, Ammonium Molybdate  |
| MOLYBDATE J POWDER      | 0.03  | Sodium Molybdate, Sulfuric Acid  |
| pH 7 BUFFER             | 1.26  |  |
| pH 10 BUFFER            | 0,1   |  |
| 2 N POTASSIUM HYDROXIDE | 0.04  |  |
| 10 N SULFURIC ACID      | 0.81  |  |
| DEIONIZED WATER         | 84.24 |  |
| TOTAL                   | 100   |  |





Philip Environmental Services Corporation 210 West Sand Bank Road Columbia, Illinois 62236

Phone: 618-281-7173 FAX: 618-281-5120

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# – MEMORANDUM –

TO: Cory Chance

FROM: Paul Anderson

DATE: January 15, 1997

PROJECT NUMBER: 17309

SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO

### SUMMARY AND DATA USABILITY

### **Data Quality Review Project Summary**

This report addresses the quality of the data results for the samples collected on September 23 and December 12, 1996, at the WFS-Milagro site.

Samples collected at the facility were analyzed for a selected list of analytical parameters which included the following: Volatiles (8260); Semi-volatiles (8270); and Metals (Ba, Cd, Cr, Pb, Ag, As, Se by 6010) and (As by 7060, Se by 7740, Hg by 7471). The analyte list and detection limits were to follow SW-846.

The analyses were performed and reported by Philip Analytical Services, Inc. (PAS), for the samples received by the laboratory on September 24 and December 13, 1996. Table 1 presents a summary of the validated samples, including their identification information, matrix, collection dates, methods of analysis performed, and deviated quality control criteria noted for each sample. The data package included a summary of sample results with QC batch summary results. Overall, the data packages were sufficiently complete to properly assess the quality of the analytical results.

### **Data Usability**

The analytical data were reviewed and qualified in accordance with U.S. Environmental Protection Agency (USEPA) Data Validation Functional Guidelines (USEPA, 1994a,b). Rejected data, which are considered unusable for either qualitative or quantitative purposes, are a result of a major deficiency noted in the data generation process. Minor deficiencies in the data generation process result in approximation of sample data. Approximation of a data point indicates uncertainty in the reported concentration of the chemical, but not its assigned identity. Conservative assumptions made when basing conclusions on analytical data allow the use of approximated analytical data. This approach to the use of analytical data is consistent 1/23/97/E/1/7309/QALABDAT.DOC

PAGE: 2 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

with USEPA (USEPA, 1992) risk assessment and data usability guidance. A summary of specific quality assurance/quality control (QA/QC) excursions that resulted in qualification of sample data are presented in the following sections.

This section summarizes the analytical data in terms of its completeness and usability for characterization of the current water quality for engineering decisions and risk management. Data completeness is defined as the percentage of sample results that have been determined to meet quality control criteria during the data review process. The completeness of the data set was determined to be 99%. However, some sample results have been qualified due to method blank, matrix spike/matrix spike duplicate (MS/MSD), and laboratory control sample (LCS) excursions described below.

Data usability is a qualitative evaluation that utilizes the findings of the data validation review, historical information, regulatory guidance and other applicable information to provide an assessment of how the data set may be used. The assessment is made with reference to the specific data uses specified in the previous paragraph. Data which do not meet completeness requirements are not considered usable, however, data which are complete are not necessarily usable.

The data usability analysis found that no issues which affect data usability were noted.

### DATA VALIDATION METHODS AND RESULTS

### **Data Validation Protocols**

The criteria employed for data validation review (USEPA, 1994) provide differing quality control criteria from those specified in the methods. In cases where method requirements and validation requirements overlap, method quality control criteria are used and the validation guidelines are followed for applying data qualifiers, if required. The data validation guidelines quality control criteria are used when the method does not provide criteria. Data may be in full compliance with the method requirements, but may not meet data validation guidance.

PAGE: 3 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

### Quality Control Parameters

The following QA/QC parameters were evaluated for organic analyses:

- 1. Holding Times and Preservation
- 2. Blank Analyses
- 3. Surrogate Spikes
- 4. MS/MSD Analysis
- 5. LCS Analysis
- 6. Reporting Limits

The following QA/QC parameters were evaluated for inorganic analyses:

- 1. Holding Times and Preservation
- 2. Blank Analyses
- 3. MS Sample Recovery Criteria
- 4. LCS Analysis
- 5. Laboratory Replicate Analysis
- 6. Reporting Limits

### Data Qualifiers Used in Validation

The following data qualifiers are defined by the USEPA in the Data Validation Functional Guidelines for use in validating analytical data. The appropriate data qualifiers were added to the data packages in red ink during data validation. Any necessary corrections to the data due to transcription errors or calculation errors were also added with red ink. Each page with changes was also initialed and dated by the validator.

The following data validation qualifiers may be used during the validation of data:

- R Indicates that the reporting limit or sample result has been determined to be unusable due to a major deficiency in the data generation process. The data should not be used for any qualitative or quantitative purposes.
- U Indicates that the analyte was analyzed for, but was not detected. The sample-specific reporting limit is presented and adjusted for dilution and percent solids, as appropriate. The qualifier is also used to signify that the reporting limit of an analyte was raised due to blank contamination.

### PAGE: 4 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

- J The analyte was positively identified but the concentration should be considered approximate. This qualifier indicates that the data validation process identified a deficiency in the data generation process.
- UJ Indicates that the sample-specific reporting limit for the analyte in this sample should be considered approximate. This qualifier is used when the data validation process identified a deficiency in the data generation process.
- N The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification".
- NJ The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

### **Data Evaluation Results**

This section describes the qualifiers that were added to sample data where QA/QC criteria were not met. QA/QC parameters that were met are not discussed. Samples that required qualification are identified in the following sections by the sample location documented on the field chain-of-custody record.

### **General Comments**

The packages included sample results along with QC results for blank analyses, matrix spikes, blank spikes, laboratory replicates, and surrogates. The packages were sufficiently complete to perform the Level I validation on the samples. The laboratory QC information was inadvertently missing from the initial package. The laboratory was notified and submitted the missing information.

The September COC 'C3115', had indicated that the metals analyses should be performed as TCLP and not as Totals. However, the laboratory supplied a faxed statement from Philip Environmental which corrected the COC request to correctly analyze the samples as Total RCRA Metals.

PAGE: 5 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

### Volatiles

### **Blank Analysis**

Contamination was detected in the method blanks. The sample results have been qualified based on the observed contamination. Sample results reported at concentrations less than five times the blank contamination (ten times for the common laboratory contaminants) have been 'U' qualified as not detected and the reporting limit raised to the reported value for the analyte. The observed compounds, their reported concentrations and the affected samples are summarized in the following table. Blank contaminants that did not affect sample results have not been listed in the table.

| Blank LD. | Analysis<br>Date | Analyte | Conc.<br>(mg/kg) | Reporting<br>Limit(mg/kg) | Action<br>Level <sup>1</sup> (mg/kg) | Affected<br>Samples                            | Qualified<br>Results <sup>2</sup>        |
|-----------|------------------|---------|------------------|---------------------------|--------------------------------------|--|--|
| 1002SM02  | 10/02/96         | acetone | 0.052            | 0.030                     | 0.520                                | BH1 5-7<br>BH1 25-27<br>BH2 17-18<br>BH3 20-22 | 0.12 U<br>0.27 U<br>0.17 U<br>0.07 U     |
|           |                  | DCM     | 0.21             | 0.010                     | 2.100                                | BH1 5-7<br>BH1 25-27<br>BH2 17-18<br>BH3 20-22 | 0.19 U<br>0.21 U<br>0.17 U<br>0.18 U     |
|           |                  | toluene | 0.006            | 0.005                     | 0.030                                | BH1 5-7<br>BH1 25-27<br>BH2 17-18              | 0.005 U<br>0.007 U<br>0.005 U            |
| 1005SM02  | 10/05/96         | acetone | 0.066            | 0.030                     | 0.660                                | BG1 5-7  | 0.072 U                                  |
|           |                  | DCM     | 0.118            | 0.010                     | 1.180                                | BG1 5-7  | 0.091 U                                  |
|           |                  | toluene | 0.006            | 0.005                     | 0.030                                | BG1 5-7  | 0.005 U                                  |
| 1220SM01  | 12/20/96         | acetone | 0.170            | 0.030                     | 1.700                                | BH1 3-5<br>BH1 10-12<br>BH2 15-17<br>BH3 13-15 | 0.18 U<br>0.15 U<br>0.14 U<br>0.12 U     |
|           |                  | DCM     | 0.020            | 0.010                     | 0.200                                | BH1 3-5<br>BH1 10-12<br>BH2 15-17<br>BH3 13-15 | 0.039 U<br>0.034 U<br>0.029 U<br>0.033 U |

Action References:

Action Level = highest blank contaminant value above the reporting limit x 5 or x 10.
 Qualified Results = reported value, which becomes the qualified reporting limit.

Notes:

#### DCM - methylene chloride, dichloromethane

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PAGE: 6 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

#### Laboratory Control Samples

Percent recoveries were outside the control limits for several of the analytes in several of the blank spike samples. For recoveries that were less than control limits, the non-detected sample results were 'UJ' qualified as estimated reporting limits. The excursions that were above the control limits did not affect the data, since there were no positive results for the excursions. LCS excursions are presented in the following table. Excursions that did not affect sample results have not been listed in the table.

| QC<br>Batch File | Analysis<br>Date | Matrix | Analyte       | Percent<br>Recovery | Control<br>Limits (%) | Action | Affected<br>Samples                                       |
|------------------|------------------|--------|---------------|---------------------|-----------------------|--------|---|
| 1002SM02         | 10/02/96         | soil   | vinyl acetate | 76                  | 80-120                | 1      | BH1 5-7<br>BH1 25-27<br>BH2 17-18<br>BH3 20-22<br>BG1 5-7 |

Action References: 1 - The non-detected results for the analyte have been 'UJ' qualified as estimated reporting limits.

Notes:

Sample BG1 5-7 has been qualified with the 1002SM02 batch data, since there was no spike data available for the 1005SM02 batch.

The lab had commented on the first narrative that the vinyl acetate low recoveries may be due to the samples sitting at room temperature for several hours, and that the compound may have degraded in that period of time. Upon calling the laboratory for more information, it was determined that this was not the case, as it sounded in the comment. The samples were simply setting in the auto-sampler tray waiting to start. Therefore, this is not a concern.

#### Matrix Spike/Matrix Spike Duplicates

MS/MSD percent recoveries were outside the control limits for several analytes for the matrix spike samples. No qualification was required for percent recoveries that exceeded control limits if there were either no positive detects for those analytes or if the positive results had a concentration that was greater than 4x the spike added. Samples were not qualified if only one of the two spikes were outside the limits. Matrix spike excursions are presented in the following table. Excursions that did not affect sample results have not been listed in the table.

PAGE: 7 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

| Sample ID | Analysis<br>Date | Matrix | Analyte       | %Recovery<br>(MS/MSD) | Control<br>Limits (%) | Action |
|-----------|------------------|--------|---------------|-----------------------|-----------------------|--------|
| BH3 13-15 | 12/20/96         | soil   | vinyl acetate | 41/28                 | 80-120                | 1      |

Action References:

1 - The non-detected result for the analyte has been 'UJ' qualified as an estimated reporting limits.

### Semi-Volatiles

Several of the samples have had their detection limits raised by a factor of two because of sample loss in the Gel Permeation Cleanup (GPC). Ten mLs of sample extract is put through the GPC. Half of the extract is lost in cleaning out the lines to avoid cross contamination. Once the sample goes through the system, it is brought back to original volume of extract, thus resulting in a two-fold dilution.

### Laboratory Control Samples

Percent recoveries were outside the control limits for several of the analytes in the blank spike samples. For recoveries that were less than control limits, the non-detected sample results were 'UJ' qualified as estimated reporting limits. The excursions that were above the control limits did not affect the data, since there were no positive results associated with the excursions. LCS excursions are presented in the following table. Excursions that did not affect sample results have not been listed in the table.

| QC<br>Batch File | Preparation<br>Date | Matrix | Analyte                | %Recovery<br>(BS/BSD) | Control<br>Limits (%) | Action | Affected<br>Samples |
|------------------|---------------------|--------|------------------------|-----------------------|-----------------------|--------|---------------------|
| 1217PB02         | 12/17/96            | soil   | 2-chlorophenol         | 19/31                 | 25-102                | 1      | BH1 3-5             |
|                  |                     |        | 1,4-dichlorobenzene    | NR / 1.0              | 28-104                | 2      | BH1 10-12           |
|                  |                     |        | 1,2,4-trichlorobenzene | 4.0 / 17              | 38-107                | 2      | BH2 15-17           |
|                  |                     |        |                        |                       |                       |        | BH3 13-15           |

Action References: 1 - The non-detected results for the analyte have been 'UJ' qualified as estimated reporting limits.

2 - The non-detected results for the analytes have been 'R' qualified as unusable.

PAGE: 8 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

### Matrix Spike/Matrix Spike Duplicates

MS/MSD percent recoveries were outside the control limits for several analytes for the matrix spike samples. No qualification was required for percent recoveries that exceeded control limits if there were either no positive detects for those analytes or if the positive results had a concentration that was greater than 4x the spike added. Samples were not qualified if only one of the two spikes were outside the limits. Matrix spike excursions are presented in the following table. Excursions that did not affect sample results have not been listed in the table.

| Sample ID | Preparation<br>Date | Matrix | Analyte                                       | %Recovery<br>(MS/MSD) | Control<br>Limits (%) | Action |
|-----------|---------------------|--------|---|-----------------------|-----------------------|--------|
| BH1 3-5   | 12/17/96            | soil   | 1,4-dichlorobenzene<br>1,2,4-trichlorobenzene | 2.0 / 3.0<br>21 / 19  | 28-104<br>38-107      | 1      |

Action References: 1 - The non-detected results for the analytes have been 'R' qualified as unusable.

### Metals

The detection limits are elevated in the samples due to a difference in the digestion for the samples and the standards. In order to match the matrices a dilution multiplier had to be applied.

Matrix Spike/Matrix Spike Duplicates

MS/MSD percent recoveries were outside the control limits for several analytes for the matrix spike samples. No qualification was required for percent recoveries that exceeded control limits if there were either no positive detects for those analytes or if the positive results had a concentration that was greater than 4x the spike added. The following table presents the matrix spike excursions. Excursions that did not affect sample results have not been listed in the table.

| Sample I.D. | QC Batch<br>Date | Matrix | Analyte             | %Recovery<br>(MS/MSD) | Control<br>Limits (%) | Action | Affected<br>Samples                                       |
|-------------|------------------|--------|---------------------|-----------------------|-----------------------|--------|---|
| BH1 5-7     | 1002MJA1         | soil   | arsenic<br>selenium | 54/71<br>70/73        | 75-125<br>75-125      | 1<br>1 | BH1 5-7<br>BH1 25-27<br>BH2 17-18<br>BH3 20-22<br>BG1 5-7 |

Action References:

1 - Detected results for the analyte have been 'J' qualified as estimated and the nondetected results for the analyte have been 'UJ' qualified as estimated reporting limits. PAGE: 9 SUBJECT: ANALYTICAL DATA QA REVIEW - WFS MILAGRO MEMO FROM: Paul Anderson DATE: January 23, 1997

### Table 1

### SAMPLE CROSS REFERENCE

| C.O.C. | Client I.D. | Lab I.D. | Package<br>I.D. | Matrix | Collection<br>Date | Methods                               | Deviated<br>Criteria     |
|--------|-------------|----------|-----------------|--------|--------------------|---------------------------------------|--------------------------|
| C3115  | BH1 5-7     | 38766    | AN961045        | soil   | 9/23/96            | Volatiles<br>Semi-volatiles<br>Metals | MB, BS<br>none<br>MS     |
| C3115  | BH1 25-27   | 38768    | AN961045        | soil   | 9/23/96            | Volatiles<br>Semi-volatiles<br>Metals | MB, BS<br>none<br>MS     |
| C3115  | BH2 17-18   | 38770    | AN961045        | soil   | 9/23/96            | Volatiles<br>Semi-volatiles<br>Metals | MB, BS<br>none<br>MS     |
| C3115  | BH3 20-22   | 38772    | AN961045        | soil   | 9/23/96            | Volatiles<br>Semi-volatiles<br>Metals | MB, BS<br>none<br>MS     |
| C3115  | BG1 5-7     | 38774    | AN961045        | soil   | 9/23/96            | Volatiles<br>Semi-volatiles<br>Metals | MB, BS<br>none<br>MS     |
| C3017  | BH1 3-5     | 55066    | AN961368        | soil   | 12/12/96           | Volatiles<br>Semi-volatiles<br>Metals | MB, MS<br>BS, MS<br>none |
| C3017  | BH1 10-12   | 55067    | AN961368        | soil   | 12/12/96           | Volatiles<br>Semi-volatiles<br>Metals | MB, MS<br>BS<br>none     |
| C3017  | BH2 15-17   | 55068    | AN961368        | soil   | 12/12/96           | Volatiles<br>Semi-volatiles<br>Metals | MB, MS<br>BS<br>none     |
| C3017  | BH3 13-15   | 55069    | AN961368        | soil   | 12/12/96           | Volatiles<br>Semi-volatiles<br>Metals | MB, MS<br>BS<br>none     |

Method References: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986, and it's updates, (SW-846 / 8260,8270, 6010, 7060, 7740, 7471).

Excursion References: MB - Method Blank

MS - Matrix Spike Recovery

BS - Blank Spike Recovery (Laboratory Control Sample)

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## Zenon Env., onmental Laboratories - Certificate of Analysis

Page 7 of 14

| Component                             | Client ID:<br>Zenon ID:<br>Date Sampled:<br>MDL | Units      | BH1-5-7<br>038766 96<br>96/09/23 | BH1<br>25-27<br>038768 96<br>96/09/23 | BH2<br>17-18<br>038770 96<br>96/09/23 |
|---------------------------------------|---|------------|----------------------------------|---------------------------------------|---------------------------------------|
| Volatiles via SW846 Method            |   | Oms        |                                  |                                       |                                       |
| Acetone                               | 0.030   | mg/kg      | 60.12 11                         | < 0.27 U                              | < 0.17 U                              |
| Acrolein                              | 0.010   | " "        | <                                | <                                     | <                                     |
| Acrylonitrile                         | 0.010   | <b>i</b> . | <                                | <                                     | <                                     |
| Benzene                               | 0.005   | 47         | <                                | <                                     | <                                     |
| Bromoform                             | 0.010   | 17         | <                                | <                                     | <                                     |
| Bromomethane                          | 0.010   | u.         | <                                | <                                     | <                                     |
| 2-Butanone                            | 0.015   | 17         | <                                | <                                     | <                                     |
| Carbon Disulfide                      | 0.010   | <b>T</b> 5 | <                                | <                                     | <                                     |
| Carbon Distantice                     | 0.010   |            | <                                | <                                     | <                                     |
| Chlorobenzene                         | 0.005   | 14         | <                                | <                                     | <                                     |
| Chlorodibromomethane                  | 0.005   | *1         | <                                | <                                     | <                                     |
| Chloroethane                          | 0.005   | a          | <                                | <                                     | <                                     |
|                                       | 0.010   |            | <                                | <                                     | <                                     |
| 2-Chloroethylvinylether<br>Chloroform | 0.005   |            | <                                | <                                     | <                                     |
| Chloromethane                         | 0.000   | н          | <                                | <                                     | <                                     |
| 1,2-Dichlorobenzene                   | 0.005   | 11         | <                                | <                                     | <                                     |
| 1,3-Dichlorobenzene                   | 0.005   | **         | <                                | <                                     | <                                     |
| 1,4-Dichlorobenzene                   | 0.005   |            | <                                | <.                                    | <                                     |
| Dichlorobromomethane                  | 0.005   | 11         | <                                | <                                     | <                                     |
| 1,1-Dichloroethane                    | 0.005   | 11         | <                                | <                                     | <                                     |
| 1,2-Dichloroethane                    | 0.005   | n          | <                                | <                                     | <                                     |
|                                       | 0.010   | 11         | <                                | <                                     | <                                     |
| 1,1-Dichloroethene                    | 0.010   |            | <                                | <                                     | <                                     |
| cis-1,2-Dichloroethene                | 0.010   | 14         | <                                | <                                     | <                                     |
| trans-1,2-Dichloroethene              | 0.005   | 17         | <                                | <                                     | <                                     |
| 1,2-Dichloropropane                   |   | 11         | <                                | <                                     | <                                     |
| cis-1,3-Dichloropropene               | 0.005   | 11         | <                                | <                                     | <                                     |
| trans-1,3-Dichloropropene             | 0.005   | *1         | <                                | <                                     | <                                     |
| Ethylbenzene                          | 0.005   |            | <                                | <                                     | <                                     |
| 2-Hexanone                            | 0.010   | ч          | <b>∠</b> 0.19 <b>U</b>           | <b>4</b> 0.21 <b>U</b>                | L0.17U                                |
| Methylene Chloride                    | 0.010   | ۰.         |                                  | <                                     | <                                     |
| 4-Methyl-2-Pentanone                  | 0.010   | **         | <                                |                                       |                                       |
| Styrene                               | 0.005   | 11         | <                                | <                                     | <                                     |
| 1,1,1,2-Tetrachloroethane             | 0.010   |            | <                                |                                       | <                                     |
| 1,1,2,2-Tetrachloroethane             | 0.010   | 11         | <                                | <<br><                                | <     <                               |
| Tetrachloroethene                     | 0.020   |            | <                                | <0.007 U                              | €0.005 U                              |
| Toluene                               | 0.005   | 11         | <b>∠</b> 0.005 <b>∪</b>          |                                       |                                       |
| 1,1,1-Trichloroethane                 | 0.005   | •1         | <                                | <                                     | <                                     |
| 1,1,2-Trichloroethane                 | 0.010   |            | <                                | <                                     | <                                     |
| Trichloroethene                       | 0.005   | 11         | <                                | <                                     | <                                     |
| Vinyl Acetate                         | 0.010   |            | < いづ                             |                                       |                                       |
| Vinyl Chloride                        | 0.010   |            | <                                | <                                     | <                                     |
| Xylenes(Total)                        | 0.005   |            | <                                | <                                     | <                                     |
| Surrogate Recoveries                  |   | 90         | 0.0                              | 0.7                                   | 00                                    |
| d4-1,2-Dichloroethane                 |   |            | 93                               | 95                                    | 99                                    |
| d8-Toluene                            |   |            | 94                               | 91                                    | 97                                    |
| Bromofluorobenzene                    |   |            | 95                               | 95                                    | 97                                    |

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# Zenon Env., onmental Laboratories - Certificate of Analysis

Page 8 of 14

|  | Client ID:    |       | BH1-5-7          | BH1<br>25-27                            | BH2<br>17-18   |
|--|---------------|-------|------------------|---|----------------|
|  | Zenon ID:     |       | 038766 96        | 038768 96                               | 038770 96      |
|  | Date Sampled: |       | 96/09/23         | 96/09/23                                | 96/09/23       |
| Component                                    | MDL           | Units | 10,01,22         | , | 10,07,20       |
| Semi - volatiles via SW846 N                 |               | Omts  |                  |   |                |
| Phenol                                       | 0.11          | mg/kg | < 0.22           | < 0.22                                  | <0.22          |
| Bis(2-chloroethyl)ether                      | 0.18          | "     | <0.36            | < 0.36                                  | < 0.36         |
| 2-Chlorophenol                               | 0.27          |       | < 0.54           | < 0.54                                  | < 0.54         |
| 1,3-Dichlorobenzene                          | 0.20          | м     | <0.40            | <0.40                                   | <0.40          |
| 1-4-Dichlorobenzene                          | 0.20          | ч     | <0.40            | <0.40                                   | <0.40          |
| 1,2-Dichlorobenzene                          | 0.20          | **    | <0.40            | <0.40                                   | <0.40          |
| Bis(2-chloroisopropyl)ether                  | 0.15          | **    | <b>&lt;0.3</b> 0 | <0.30                                   | < 0.30         |
| Hexachloroethane                             | 0.20          | ч     | <0.50            | <0.40                                   | < 0.40         |
| N-Nitroso-di-N-Propylamine                   | 0.20          |       | <0.40            | <0.42                                   | <0.40          |
| Nitrobenzene                                 | 0.20          |       | <0.42<br><0.40   | <0.42                                   | <0.42          |
| Isophorone                                   | 0,40          | a     | <0.40            | <0.40                                   | <0.40<br><0.80 |
| 2-Nitrophenol                                | 0,14          | J.r   | <0.80            | <0.80                                   | <0.80          |
| 2.4-Dimethylphenol                           | 0.17          | 11    | < 0.34           | < 0.28                                  | < 0.28         |
| Bis(2-chloroethoxy)methane                   | 0.13          | 11    | <0.26            | <0.26                                   | <0.26          |
| 2,4-Dichlorophenol                           | 0.13          | IT    | <0.20            | <0.20<br><0.24                          | <0.20          |
| 1,2,4-Trichlorobenzene                       | 0.12          | 11    | <0.24<br><0.40   | < 0.40                                  | <0.24          |
| Naphthalene                                  | 0.03          |       | <0.40<br><0.06   | <0.40                                   | <0.40          |
| Hexachlorobutadiene                          | 0.03          | .,    | <0.00<br><0.40   | <0.00<br><0.40                          | <0.00          |
| 4-Chloro-3-Methylphenol                      | 0.14          | ,,    | <0.40            | <0.28                                   | <0.28          |
|  |               | ,1    | <0.28            | < 0.26                                  | <0.26          |
| Hexachlorocyclopentadiene                    | 0.20          | 81    |                  | <0.40<br><0.24                          | <0.40          |
| 2,4,6-Trichlorophenol<br>2-Chloronaphthalene | 0.12<br>0.09  | *1    | <0.24<br><0.18   | < 0.18                                  | <0.24          |
| Acenaphthylene                               | 0.09          |       | <0.18            | <0.18                                   | <0.18          |
| Dimethyl phthalate                           | 0.11          | .,    | <0.22            | <0.08                                   | <0.22          |
| 2,6-Dinitrotoluene                           | 0.06          | **    | <0.12            | <0.12                                   | <0.22          |
| Acenaphthene                                 | 0.00          |       | <0.12            | <0.12                                   | <0.12          |
| 2,4-Dinitrophenol                            | 0.07          |       | <0.96            | <0.14                                   | <0.14          |
| 2,4-Dinitrotoluene                           | 0.48          | 11    | <0.90<br><0.10   | <0.10                                   | <0.90          |
| 4-Nitrophenol                                |               | 11    | <0.10            | <0.10                                   | <0.10          |
| Fluorene                                     | 0.14          | "     | <0.26<br><0.06   | <0.28                                   |                |
|  | 0.03          | н     |                  |   | < 0.06         |
| 4-Chlorophenylphenylether                    | 0.09          | 14    | <0.18            | < 0.18                                  | <0.18          |
| Diethyl phthalate                            | 0.11          |       | <0.22            | < 0.22                                  | <0.22          |
| 4,6-Dinitro-2-methylphenol                   | 0.15          |       | < 0.30           | <0.30                                   | <0.30          |
| N-Nitrosodiphenylamine                       | 0.19          | 13    | <0.38            | <0.38                                   | <0.38          |
| 4-Bromophenylphenylether                     | 0.03          | 17    | <0.06            | <0.06                                   | <0.06          |
| Hexachlorobenzene                            | 0.20          | 1     | <0.40            | <0.40                                   | <0.40          |
| Pentachlorophenol                            | 0.11          | 11    | < 0.22           | < 0.22                                  | <0.22          |
| Phenanthrene                                 | 0.03          |       | <0.06            | <0.06                                   | <0.06          |
| Anthracene                                   | 0.02          |       | <0.04            | < 0.04                                  | < 0.04         |
| Di-n-butyl phthalate                         | 0.11          |       | <0.22            | < 0.22                                  | < 0.22         |
| Fluoranthene                                 | 0.02          |       | < 0.04           | < 0.04                                  | < 0.04         |
| Pyrene                                       | 0.03          |       | <0.06            | < 0.06                                  | <0.06          |
| Benzyl butyl phthalate                       | 0.06          |       | < 0.12           | < 0.12                                  | < 0.12         |
| 3,3-Dichlorobenzidine                        | 0.10          |       | <0.20            | <0.20                                   | < 0.20         |
| Benzo(a)anthracene                           | 0.02          |       | <0.04            | <0.04                                   | < 0.04         |
| Chrysene                                     | 0.03          |       | <0.06            | <0.06                                   | <0.06          |

Page 9 of 14

10/9/96

## Zenon Env., onmental Laboratories - Certificate of Analysis

|                            |               |       |           | BH1       | BH2       |
|----------------------------|---------------|-------|-----------|-----------|-----------|
|                            | Client ID:    |       | BH1-5-7   | 25-27     | 17-18     |
|                            | Zenon ID:     |       | 038766 96 | 038768 96 | 038770 96 |
|                            | Date Sampled: |       | 96/09/23  | 96/09/23  | 96/09/23  |
| Component                  | MDL           | Units |           |           |           |
| Bis(2-ethylhexyl)phthalate | 0.14          | ••    | <0.28     | <0.28     | <0.28     |
| Di-n-octyl phthalate       | 0.11          | •     | <0.22     | <0.22     | < 0.22    |
| Benzo(b)fluoranthene       | 0.04          | "     | <0.08     | <0.08     | <0.08     |
| Beuzo(k)fluoranthene       | 0.04          | **    | <0.08     | <0.08     | <0.08     |
| Benzo(a)pyrene             | 0.05          | n     | <0.10     | <0.10     | <0.10     |
| Indeno(1,2,3-cd)pyrene     | 0.06          | н     | < 0.12    | <0.12     | < 0.12    |
| Dibenzo(a,h)anthracene     | 0.04          | ч     | <0.08     | <0.08     | <0.08     |
| Benzo(ghi)perylene         | 0.04          | 11    | <0.08     | < 0.08    | <0.08     |
| N-Nitrosodimethylamine     | 1.0           | 11    | <2.0      | <2.0      | <2.0      |
| Aniline                    | 0.50          | н     | <1.0      | <1.0      | <1.0      |
| Benzyl alcohol             | 0.50          | "     | <1.0      | <1.0      | <1.0      |
| Carbazole                  | 0.50          | и     | <1.0      | <1.0      | <1.0      |
| 2-Methylphenol             | 0.50          | 9     | <1.0      | <1.0      | <1.0      |
| Benzoic acid               | 0.50          | ۰r    | <1.0      | <1.0      | <1.0      |
| 4-Chloroaniline            | 0.50          | 14    | <1.0      | <1.0      | <1.()     |
| 2-Methylnaphthalene        | 0.10          | н     | < 0.20    | < 0.20    | <0.20     |
| 2,4,5-Trichlorophenol      | 0.10          | 14    | <0.20     | <0.20     | <0.20     |
| 2-Nitroaniline             | 0.50          | "     | <1.0      | <1.0      | <1.0      |
| 3-Nitroaniline             | 0.50          | 11    | <1.0      | <1.0      | <1.0      |
| Dibenzofuran               | 0.50          | 44    | <1.0      | <1.0      | <1.0      |
| Benzidine                  | 0.50          |       | <1.0      | <1.0      | <1.0      |
| 4-Nitroaniline             | 0.50          |       | <1.0      | <1.0      | <1.0      |
| Surrogate Recoveries       |               | %     |           |           |           |
| 2-Fluorophenol             |               |       | 47        | 73        | 52        |
| d5-Phenol                  |               |       | 66        | 107       | 73        |
| d5-Nitrobenzene            |               |       | 60        | 100       | 60        |
| 2-Fluorobiphenyl           |               |       | 65        | 101       | 79        |
| 2,4,6-Tribromophenol       |               |       | 69        | 103       | 72        |
| d14-p-Terphenyl            |               |       | 83        | 128       | 82        |

Client:Philip Environmental Inc. Project:16766

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Zenon Env., onmental Laboratories - Certificate of Analysis

Page 10 of 14

|                                 |            |       |                                  | 1            |
|---------------------------------|------------|-------|----------------------------------|--------------|
|                                 |            |       | BH3                              | $\mathbf{V}$ |
|                                 | Client ID: |       | 20-22                            | BG1-5-7      |
| 2                               | Lenon ID:  |       | 038772 96                        | 038774 96    |
|                                 | Sampled:   |       | 96/09/23                         | 96/09/23     |
| Component                       | MDL        | Units | 50,00725                         | 10/07/20     |
| Volatiles via SW846 Method 8260 | MP L       | Onus  |                                  |              |
| Acetone                         | 0.030      | mg/kg | < 0.070 U                        | ∠ 0.072 U 🔆  |
| Acrolein                        | 0.010      | "     | <                                | <            |
| Acrylonitrile                   | 0.010      | л     | <                                | <            |
| Benzene                         | 0.005      | 11    | <                                | <            |
| Bromoform                       | 0.010      | ч     | <                                | <            |
| Bromomethane                    | 0.010      |       | <                                | <            |
| 2-Butanone                      | 0.015      |       | <                                | <            |
| Carbon Disulfide                | 0.010      |       | <                                | <            |
| Carbon Tetrachloride            | 0.010      | п     | <                                | <            |
| Chlorobenzene                   | 0.005      | "     | <                                | <            |
| Chlorodibromomethane            | 0.005      | 11    | <                                | <            |
| Chloroethane                    | 0.010      | 11    | <                                | <            |
| 2-Chloroethylvinylether         | 0.010      |       | <                                | <            |
| Chloroform                      | 0.005      | 1:    | <                                | <            |
| Chloromethane                   | 0.010      | н     | <                                | <            |
| 1,2-Dichlorobenzene             | 0.005      | в     | <                                | <            |
| 1,3-Dichlorobenzene             | 0.005      | 11    | <                                | <            |
| 1,4-Dichlorobenzene             | 0.005      | 14    | <                                | 0.008 🖌      |
| Dichlorobromomethane            | 0.005      | 11    | <                                | < 7          |
| 1,1-Dichloroethane              | 0.005      |       | <                                | <            |
| 1,2-Dichloroethane              | 0.005      | D.    | <                                | <            |
| 1,1-Dichloroethene              | 0.010      |       | <                                | <            |
| cis-1,2-Dichloroethene          | 0.010      |       | <                                | <            |
| trans-1,2-Dichloroethene        | 0.010      | 73    | <                                | <            |
| 1.2-Dichloropropane             | 0.005      | и     | <                                | <            |
| cis-1,3-Dichloropropene         | 0.005      | 11    | <                                | <            |
| trans-1,3-Dichloropropene       | 0.005      |       | <                                | <            |
| Ethylbenzene                    | 0.005      |       | <                                | <            |
| 2-Hexanone                      | 0.010      | 4     | <                                | 4            |
| Methylene Chloride              | 0.010      | •,    | 2 0.18 V                         | ≤ 0.091 VX   |
| 4-Methyl-2-Pentanone            | 0.010      |       | <                                | <            |
| Styrene                         | 0.005      | •     | <                                | <            |
| 1,1,1,2-Tetrachloroethane       | 0.010      | u     | <                                | <            |
| 1,1,2,2-Tetrachloroethane       | 0.010      |       | <                                | <            |
| Tetrachloroethene               | 0.020      | 11    | <                                | < ,          |
| Toluene                         | 0.005      |       | <                                | ∠ 0.005 U ¥  |
| 1,1,1-Trichloroethane           | 0.005      |       | <                                | <            |
| 1,1,2-Trichloroethane           | 0.010      | н     | <                                | <            |
| Trichloroethene                 | 0.005      | и     | <                                | <            |
| Vinyl Acetate                   | 0.010      | 11    | <v5< td=""><td>&lt;05</td></v5<> | <05          |
| Vinyl Chloride                  | 0.010      | n     | <                                | <            |
| Xylenes(Total)                  | 0.005      | **    | <                                | 0.005        |
| Surrogate Recoveries            |            | %     |                                  | •            |
| d4-1,2-Dichloroethane           |            |       | 103                              | 99           |
| d8-Toluene                      |            |       | 103                              | 100          |
| Bromofluorobenzene              |            |       | 100                              | 96           |
|                                 |            |       |                                  |              |

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# Zenon Env., onmental Laboratories - Certificate of Analysis

Page 11 of 14

|   | Client ID:    |       | BH3<br>20-22     | BG1-5-7        |
|---|---------------|-------|------------------|----------------|
|   | Zenon ID:     |       | 038772 96        | 038774 96      |
|   | Date Sampled: |       | 96/09/23         | 96/09/23       |
| Component                                 | MDL           | Units |                  |                |
| Semi - volatiles via SW846 N              |               | 0     |                  |                |
| Phenol                                    | 0.11          | mg/kg | < 0.22           | < 0.22         |
| Bis(2-chloroethyl)ether                   | 0.18          | н     | < 0.36           | < 0.36         |
| 2-Chlorophenol                            | 0.27          | 11    | < 0.54           | < 0.54         |
| 1,3-Dichlorobenzene                       | 0.20          | **    | <0.40            | <0.40          |
| 1-4-Dichlorobenzene                       | 0.20          | n     | <0.40            | <0.40          |
| 1,2-Dichlorobenzene                       | 0.20          | "     | <0.40            | <0.40          |
| Bis(2-chloroisopropyl)ether               | 0.15          | 11    | <0.30            | <0.30          |
| Hexachloroethane                          | 0.20          | "     | <0.40            | <0.40          |
| N-Nitroso-di-N-Propylamine                | 0.21          | n     | <0.42            | <0.42          |
| Nitrobenzene                              | 0.20          | 11    | <0.40            | <0.40          |
| Isophorone                                | 0.40          | -11   | <0.80            | <0.80          |
| 2-Nitrophenol                             | 0.14          |       | <0.28            | <0.28          |
| 2.4-Dimethylphenol                        | 0.17          | "     | < 0.34           | < 0.34         |
| Bis(2-chloroethoxy)methane                | 0.13          |       | <0.26            | < 0.26         |
| 2,4-Dichlorophenol                        | 0.12          |       | <0.24            | <0.24          |
| 1,2,4-Trichlorobenzene                    | 0.20          | ••    | <0.40            | <0.40          |
| Naphthalene                               | 0.03          | **    | <0.06            | < 0.06         |
| Hexachlorobutadiene                       | 0.20          | 42    | <b>&lt;0.4</b> 0 | <0.40          |
| 4-Chloro-3-Methylphenol                   | 0.14          | "     | <0.28            | < 0.28         |
| Hexachlorocyclopentadiene                 | 0.20          | *1    | <0.40            | <0.40          |
| 2,4,6-Trichlorophenol                     | 0.12          | 11    | <0.24            | <0.24          |
| 2-Chloronaphthalene                       | 0.09          |       | <0.18            | <0.18          |
| Acenaphthylene                            | 0.04          |       | <0.08            | <0.08          |
| Dimethyl phthalate                        | 0.11          | •1    | <0.22            | <0.22          |
| 2,6-Dinitrotoluene                        | 0.06          | н     | <0.12            | <0.12          |
| Acenaphthene                              | 0.07          | 11    | <0.14            | <0.14          |
| 2,4-Dinitrophenol                         | 0.48          | 4     | <0.96            | <0.96          |
| 2,4-Dinitrotoluene                        | 0.05          | u.    | <0.10            | <0.10          |
| 4-Nitrophenol                             | 0,14          |       | <0.28            | <0.28          |
| Fluorene                                  | 0.03          | н     | <0.06            | <0.06          |
| 4-Chlorophenylphenylether                 | 0.09          |       | <0.18            | < 0.18         |
| Diethyl phthalate                         | 0.11          | "     | <0.22            | < 0.22         |
| 4,6-Dinitro-2-methylphenol                | 0.15          | 11    | <0.30            | < 0.30         |
| N-Nitrosodiphenylamine                    | 0.19          | "     | <0.38            | < 0.38         |
| 4-Bromophenylphenylether                  | 0.03          | "     | <0.06            | <0.06          |
| Hexachlorobenzene                         | 0.20          | 11    | <0.40            | <0.40          |
| Pentachlorophenol                         | 0.11          |       | < 0.22           | < 0.22         |
| Phenanthrene                              | 0.03          |       | <0.06            | <0.06          |
| Anthracene                                | 0.02          |       | <0.04            | < 0.04         |
| Di-n-butyl phthalate                      | 0.11          |       | <0.22            | <0.22          |
| Fluoranthene                              | 0.02          |       | <0.04            | <0.04          |
| Pyrene<br>Deserved between the back shows | 0.03          | n     | <0.06<br><0.12   | <0.06          |
| Benzyl butyl phthalate                    | 0.06          |       | <0.12            | <0.12          |
| 3,3-Dichlorobenzidine                     | 0.10          |       | <0.20<br><0.04   | <0.20<br><0.04 |
| Benzo(a)anthracene                        | 0.02          | "     | <0.04<br><0.06   | <0.04<br><0.06 |
| Chrysene                                  | 0.03          |       | <b>N.</b> 00     | <b>NU.UU</b>   |

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# Zenon Env., onmental Laboratories - Certificate of Analysis

Page 12 of 14

|                            |               |       | BH3       |           |
|----------------------------|---------------|-------|-----------|-----------|
|                            | Client ID:    |       | 20-22     | BG1-5-7   |
|                            | Zenon ID:     |       | 038772 96 | 038774 96 |
|                            | Date Sampled: |       | 96/09/23  | 96/09/23  |
| Component                  | MDL           | Units |           |           |
| Bis(2-ethylhexyl)phthalate | 0.14          | ч     | <0.28     | <0.28     |
| Di-n-octyl phthalate       | 0.11          | •1    | < 0.22    | <0.22     |
| Benzo(b)fluoranthene       | 0.04          | **    | < 0.08    | <0.08     |
| Benzo(k)fluoranthene       | 0.04          | .,    | <0.08     | <0.08     |
| Benzo(a)pyrene             | 0.05          | н     | <0.10     | <0.10     |
| Indeno(1,2,3-cd)pyrene     | 0.06          | 11    | < 0.12    | <0.12     |
| Dibenzo(a,h)anthracene     | 0.04          | u     | <0.08     | <0.08     |
| Benzo(ghi)perylene         | 0.04          | 11    | <0.08     | <0.08     |
| N-Nitrosodimethylamine     | 1.0           | v     | <2.0      | <2.0      |
| Aniline                    | 0.50          | п     | <1.0      | <1.0      |
| Benzyl alcohol             | 0.50          |       | <1.0      | <1.0      |
| Carbazole                  | 0.50          | • 1   | <1.0      | <1.0      |
| 2-Methylphenol             | 0.50          | н     | <1.0      | <1.0      |
| Benzoic acid               | 0.50          |       | <1.0      | <1.0      |
| 4-Chloroaniline            | 0.50          | 11    | <1.0      | <1.0      |
| 2-Methylnaphthalene        | 0.10          | "     | <0.20     | <0.20     |
| 2,4,5-Trichlorophenol      | 0.10          |       | <0.20     | <0.20     |
| 2-Nitroaniline             | 0.50          | 11    | <1.0      | <1.0      |
| 3-Nitroaniline             | 0.50          | 11    | <1.0      | <1.0      |
| Dibenzofuran               | 0.50          | "     | <1.0      | <1.0      |
| Benzidine                  | 0.50          | 14    | <1.0      | <1.0      |
| 4-Nitroaniline             | 0.50          | •     | <1.0      | <1.0      |
| Surrogate Recoveries       |               | %     |           |           |
| 2-Fluorophenol             |               |       | 47        | 53        |
| d5-Phenol                  |               |       | 67        | 72        |
| d5-Nitrobenzene            |               |       | 59        | 67        |
| 2-Fluorobiphenyl           |               |       | 65        | 71        |
| 2,4,6-Tribromophenol       |               |       | 68        | 62        |
| d14-p-Terphenyl            |               |       | 82        | 82        |

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## L Summary of Analysis Pre. Daws

Page MS-13 of 14

| Batch Code:     | 0925SPA1  | 0925SPA1  | 0925SPA1     | 09258PA 1      | 0925SPA1  |
|-----------------|-----------|-----------|--------------|----------------|-----------|
| pH              | 038766 96 | 038768 96 | 038770 96    | 038772 96      | 038774 96 |
| Date analyzed   | 96/09/26  | 96/09/26  | 96/09/26     | 96/09/26       | 96/09/26  |
| Date prepared   | 96/09/25  | 96/09/25  | 96/09/25     | 96/09/25       | 96/09/25  |
| Batch Code:     | 1002MJA1  |           |              |                |           |
| Arsenic (gfaa)  | 038764 96 |           |              |                |           |
|                 | 038766 96 |           |              |                |           |
|                 | 038768 96 |           |              |                |           |
|                 | 038770 96 |           |              |                |           |
|                 | 038772 96 |           |              |                |           |
|                 | 038774 96 |           |              |                |           |
| Date analyzed   | 96/10/03  |           |              |                |           |
| Date prepared   | 96/10/02  |           |              |                |           |
| Batch Code:     | 1001MGA1  |           |              |                |           |
| Мегсшу          | 038764 96 |           |              |                |           |
| -               | 038766 96 |           |              |                |           |
|                 | 038768 96 |           |              |                |           |
|                 | 038770 96 |           |              |                |           |
|                 | 038772 96 |           |              |                |           |
|                 | 038774 96 |           |              |                |           |
| Date analyzed   | 96/10/02  |           |              |                |           |
| Date prepared   | 96/10/01  |           |              |                |           |
| Batch Code:     | 1002MJA1  |           |              |                |           |
| Selenium (gfaa) | 038764 96 |           |              |                |           |
|                 | 038766 96 |           |              |                |           |
|                 | 038768 96 |           |              |                |           |
|                 | 038770 96 |           |              |                |           |
|                 | 038772 96 |           |              |                |           |
|                 | 038774 96 |           |              |                |           |
| Date analyzed   | 96/10/04  |           |              |                |           |
| Date prepared   | 96/10/02  |           |              |                |           |
| Batch Code:     | 1002MJA1  |           |              |                |           |
| Metals          | 038764 96 |           |              |                |           |
|                 | 038766 96 |           |              |                |           |
|                 | 038768 96 |           |              |                |           |
|                 | 038770 96 |           |              |                |           |
|                 | 038772 96 |           |              | reh            |           |
|                 | 038774 96 |           |              | the lot        |           |
| Date analyzed   | 96/10/02  |           |              | ت بالصحيح المل |           |
| Date prepared   | 96/10/02  |           | O blank 10   | 02 CM 1-10     |           |
| Batch Code:     | 1002SM02  | 1005SM02  | - history -  | 1 1 1          | <u>^</u>  |
| Volatiles       | 038764 96 | 038774 96 | history wear | Mine I And     | ζ.        |
|                 | 038766 96 |           | ۲            | ما بسور        |           |
|                 | 038768 96 |           |              | 120            |           |

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|------|-----|----|----|----|

|                | 038770 96 |          |
|----------------|-----------|----------|
|                | 038772 96 |          |
| Date analyzed  | 96/10/02  | 96/10/05 |
| Date prepared  | 96/10/02  | 96/10/05 |
| Batch Code:    | 0925PB01  |          |
| Semi-volatiles | 038764 96 |          |
|                | 038766 96 |          |
|                | 038768 96 |          |
|                | 038770 96 |          |
|                | 038772 96 |          |
|                | 038774 96 |          |
| Date analyzed  | 96/09/27  |          |
| Date prepared  | 96/09/25  |          |
|                |           |          |

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5555 North Service Road Burlington, Ontario, Canada L7L 5H7 Tel: (905) 332-8788 Fax: (905) 332-9169

## Certificate of Analysis

### CLIENT INFORMATION

### LABORATORY INFORMATION

| Attention:<br>Client Name:<br>Project:<br>Project Desc: | Cory Chance<br>Philip Environmental Inc.<br>16766<br>WFS Milagro | Contact:<br>Project:<br>Date Received:<br>Date Reported: | Ada Blythe, B.Sc., C.Chem.<br>AN961045<br>96/09/24<br>96/10/09 |
|---|--|--|--|
| Address:  | 4000 Monroe Road<br>Farmington, NM<br>87401                      | Submission No.:<br>Sample No.:                           | 610684<br>038764-038774  |
| Fax Number:   | 505 326-2388   |  |  |
| Phone Number:   | 505 326-2262   |  |  |
| NOTES:  | "." = not unalysed '<' = less than                               | Method Detection Limit                                   | (MDL) 'NA' = no data available                                 |
|   | 100 can by determined for all and                                |  |  |

I.OQ can by determined for all analytes by multiplying the appropriate MDL X 3,33 Solids data is based on dry weight except for biota analyses. Organic analyses are not corrected for extruction recovery standards except for isotope dilution methods, (i.e. CARB 439 PAH, all PCDD/F and DBD/DBF analyses)

Methods used by Zenon are based upon those found in 'Standard Methods for the Examination of Water and Wastewater', Seventeenth Edition. Other methods are based on the principles of MISA or EPA methodologies.

All work recorded herein has been done in accordance with normal professional standards using accepted testing methodologies, quality assurance and quality control procedures except where otherwise agreed to by the client and testing company in writing. Any and all use of these test results shall be limited to the actual cost of the pertinent analysis done. There is no other warranty expressed or implied. Your samples will be retained at Zenon for a period of three weeks from receipt of data or as per contract.

#### COMMENTS:

"NR" Not recovered. Vinyl acetate recoveries are low, likely due to the fact that the samples sat at room temperature for several hours and the compound had degraded. "\*"Matrix interference suspected

Some blank spike recoveries are high for volatiles. We are checking the reference solution. This will have no impact on the volatiles sample data.

For Lead and Chromium, the detection limits were raised by a factor of two in the samples. This occurs because the samples need to be "matrix matched" to the standards. The digested sample is in a 20% Nitric acid matrix, whereas the standards are in 10% Nitric acid. Therefore, the sample is brought to a final concentration of 10% Nitric acid (via a two fold dilution) resulting in a two-fold increase in the detection limit.

Certified by

P42,2597 Page 1

A division of PHILIP Analytical Services Corporation

Additional comments regarding samples 038764-038774

All of the samples analysed for semi-volatiles were cleaned up using Gel Permeation Chromatography (GPC). The nature of GPC clean-up involves taking an extract at 10 mL final volume and injecting this extract into the GPC to clean out any interferences in the sample by employing the use of gel permeation chromatography. Half of that extract is lost in the GPC (ie. 5mLs) to flush the system and avoid cross contamination from the previous sample. The final volume after the GPC is performed is the same as the initial volume (ie. 10 mLs) due to dilution from solvents in the GPC procedure. Therefore, half of the extract is taken to the same initial volume (ie. 10 mLs) which results in a two-fold increase in the detection limit

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Page 2 of 14

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| Component                       | Client ID:<br>Zenon ID:<br>Date Sampled:<br>MDL | Units    | Method<br>Blank<br>038764 96<br>96/09/23 | Blank<br>Spike 1<br>038764 96<br>96/09/23 | Blank<br>Spike 1<br>038764 96<br>96/09/23<br>% Rccovery | BH1-5-7<br>038766 96<br>96/09/23 | BH1-5-7<br>038766 96<br>96/09/23<br>Duplicate | BH1-5-7<br>038766 96<br>96/09/23<br>M. Spike | BH1-5-7<br>038766 96<br>96/09/23<br>MS % Rec. |
|---------------------------------|---|----------|--|---|---|----------------------------------|---|--|---|
| Arcenic (vfaa)via Method 7060   | 0.5   | mg/kg    | v  | 4.1                                       | 83  | 1.4 J                            | ۲.4<br>۲                                      |  | 54*   |
| Mercury via Method 7471         | 0.04  | ;<br>; ≠ | v  | 0.98                                      | 98  | v                                |   | 0.98   | 98  |
| Selenium (gfaa) via Method 7740 | 0.5   | =        | V  | 4.4                                       | 88  | よっ、                              |   |  | Ð   |
| Metals via SW846 Method 6010    |   |          |  |   |   |                                  |   |  |   |
| Barium                          | 0.1   | ing/kg   | v  | 100                                       | 100   | 74                               | 99  | 270  | 100   |
| Cadmium                         | 0.2   | )<br>; = |  | 50  | 100   | 0.4                              | <0.4  | 100  | 100   |
| Chromium                        | i.  | ÷        | v  | 110                                       | 110   | <10                              | <10   | 210  | 110   |
| T ead                           | 10  | =        | v  | 100                                       | 100   | <20                              | <20   | 210  | 110   |
| Silver                          | 0.5   | ÷        | v  | 52  | 100   | <1.0                             | <1.0  | 100  | 100   |

Client: Philip Environmental Inc. Project: 16766

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| Page 3 of 14   |  |  |
|--|--|--|
| Page   |  |  |
|  | BG1-5-7<br>038774 96<br>96/09/23<br>(1)<br>2.2 J<br>< v<br>>   | 140<br><0.4<br><10<br><1.0<br><1.0   |
| alysis   | BH3<br>20-22<br>038772 96<br>96/09/23<br>(1)<br>1.2 J<br>< v J   | 160<br>0.4<br><10<br><20<br><1.0   |
| ate of An  | BH2<br>17-18<br>038770 96<br>96/09/23<br>(1)<br>1.2 T<br>< u J   | 110<br><0.4<br><10<br><20<br><1.0  |
| s - Certific   | BH1<br>25-27<br>038768 96<br>96/09/23<br>(1)<br>1.3 J<br>< u J   | 96<br>0.5<br><20<br><1.0   |
| Zenon Environmental Laboratories - Certificate of Analysis | BH1-5-7<br>038766 96<br>96/09/23<br>MSD % Rec.   | 98<br>100<br>98  |
| nental La  | BH1-5-7<br>038766 96<br>96/09/23<br>MS Dup<br>5.0<br>0.99<br>3.8   | 270<br>98<br>210<br>98   |
| ıvironı  | Units<br>mg/kg<br>"  | д<br>20/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2<br>2/2<br>2/2<br>2/2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2/2<br>2<br>2<br>2/2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2<br>2 |
| Zenon Ei   | Client ID:<br>Zenon ID:<br>Date Sampled:<br>MDL<br>0.5<br>0.04<br>0.5                                    | 0.1<br>5<br>0.5<br>0.5   |
| 96/6/01  | Component<br>Arscnic (gfaa)via Method 7060<br>Mercury via Method 7471<br>Selenium (gfaa) via Method 7740 | Metals via SW846 Method 6010<br>Barium<br>Cadmium<br>Chromium<br>Lead<br>Silver  |

Client: Philip Environmental Inc. Project: 16766

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# Zenon Env. unmental Laboratories - Certificate of Analysis

Page 4 of 14

|                           | Client ID:<br>Zenon ID:<br>Date Sampled: |       | Method<br>Blank<br>038764 96<br>96/09/23 | Blank<br>Spike 1<br>038764 96<br>96/09/23 | Blank<br>Spike 1<br>038764 96<br>96/09/23 | Blank<br>Spike 2<br>038764 96<br>96/09/23 | Blank<br>Spike 2<br>038764 96<br>96/09/23 |
|---------------------------|--|-------|--|---|---|---|---|
| Component                 | MDL                                      | Units | of.                                      | 96/09/23                                  | % Recovery                                |   | % Recovery                                |
| Volatiles via SW846 Metho | d 8260                                   |       | 10                                       | (•  |   |   |   |
| Acetone                   | 0.030                                    | mg/kg | (0.0526 0                                | 0.50                                      | (170)                                     | -   | -   |
| Acrolein                  | 0.010                                    | 11    | $\sim$                                   | 0.93                                      | (190)                                     | -   | -   |
| Acrylonitrile             | 0.010                                    | u.    | <  | 0.56                                      | 110                                       | -   | -   |
| Benzene                   | 0.005                                    | +1    | <  | 0.24                                      | 98  | -   | -   |
| Bromoform                 | 0.010                                    | 11    | <  | 0.24                                      | 94  | -   | -   |
| Bromomethane              | 0.010                                    | 11    | <  | 0.25                                      | 100                                       | -   | -   |
| 2-Butanone                | 0.015                                    | 11    | <  | 0.47                                      | (190)                                     | -   |   |
| Carbon Disulfide          | 0.010                                    | 11    | <  | 0.35                                      | (140                                      | _   | _   |
| Carbon Tetrachloride      | 0.010                                    | 11    | <  | 0.23                                      | 93  | -   | -   |
| Chlorobenzene             | 0.005                                    | ,,    | <  | 0.23                                      | 92  | -   | _   |
| Chlorodibromomethane      | 0.005                                    | 11    | <  | 0.24                                      | 96  | -   | _   |
| Chloroethane              | 0.010                                    | 11    | <  | 0.25                                      | 98  | 2   | -   |
| 2-Chloroethylvinylether   | 0.010                                    | а     | <  | 0.24                                      | 97  | _   | _   |
| Chloroform                | 0.005                                    | u     | <  | 0.24                                      | 98  | -   | -   |
| Chloromethane             | 0.010                                    | 11    | <  | 0.27                                      | 110                                       | -   | -   |
| 1,2-Dichlorobenzene       | 0.005                                    | н     | ~  | 0.21                                      | 84  | _   | _   |
| 1,3-Dichlorobenzene       | 0.005                                    |       | <  | 0.21                                      | 85  | -   | -   |
| 1,4-Dichlorobenzene       | 0.005                                    |       | <  | 0.21                                      | 83  | -   | -   |
| Dichlorobromomethane      | 0.005                                    |       | <  | 0.24                                      | 84<br>94                                  | -   | -   |
| 1,1-Dichloroethane        | 0.005                                    |       | <  | 0.24                                      | 99  | -   | -   |
| 1.2-Dichloroethane        | 0.005                                    |       |  | 0.23                                      | 99<br>92                                  | -   | -   |
| 1,1-Dichloroethene        | 0.010                                    | п     | <  | 0.23                                      | (150)                                     | -   | -   |
| cis-1,2-Dichloroethene    |  | "     | <  |   |   | -   | -   |
| trans-1,2-Dichloroethene  | 0.010                                    | 11    | <  | 0.24                                      | 96  | -   | -   |
|                           | 0.010                                    | 11    | <  | 0.35                                      | 140                                       | -   | -   |
| 1,2-Dichloropropane       | 0.005                                    |       | <  | 0.23                                      | 94  | -   | -   |
| cis-1,3-Dichloropropene   | 0.005                                    | 0     | <  | 0.28                                      | 110                                       | -   | -   |
| trans-1,3-Dichloropropene | 0.005                                    | 47    | <  | 0.25                                      | 98  | -   | -   |
| Ethylbenzene              | 0.005                                    |       | <  | 0.24                                      | 95  | -   | -   |
| 2-Hexanone                | 0.010                                    | **    | < 19                                     | 0.32                                      | 030                                       | -   | -   |
| Methylene Chloride        | 0.010                                    | ч     | (0.21) <b>0</b> , 1                      | 0.53                                      | 120                                       | -   | -   |
| 4-Methyl-2-Pentanone      | 0.010                                    | *     | <  | 0.22                                      | 89  | -   | -   |
| Styrene                   | 0.005                                    | **    | <  | 0.24                                      | 97  | ~   | -   |
| 1,1,1,2-Tetrachloroethane | 0.010                                    | 11    | <  | 0.23                                      | 90  | -   | -   |
| 1,1,2,2-Tetrachloroethane | 0.010                                    |       | <  | 0.21                                      | 82  | -   | -   |
| Tetrachloroethene         | 0.020                                    | ••    |  | 0.21                                      | 86  | -   | -   |
| Toluene                   | 0.005                                    |       | 0.006                                    | 0.24                                      | 97  | -   | -   |
| 1,1,1-Trichloroethane     | 0.005                                    | "     |  | 0.23                                      | 93  | -   | -   |
| 1,1,2-Trichloroethane     | 0.010                                    | **    | <  | 0.23                                      | 94  | -   | -   |
| Trichloroethene           | 0.005                                    | н     | <  | 0.24                                      | 95  | -   | -   |
| Vinyl Acetate             | 0.010                                    | *1    | <  | 0.19                                      | (76)                                      | -   | -   |
| Vinyl Chloride            | 0.010                                    | 14    | <  | 0.29                                      | ΠÓ  | -   | -   |
| Xylenes(Total)            | 0.005                                    | "     | <  | 0.50                                      | 100                                       | -   |   |
| Surrogate Recoveries      |  | %     |  |   |   |   |   |
| d4-1,2-Dichloroethane     |  |       | 93                                       | 102                                       | 102                                       | -   | -   |
| d8-Toluene                |  |       | 98                                       | 100                                       | 100                                       | -   | - /                                       |
| Bromofluorobenzene        |  |       | 96                                       | 98  | 98  | -   | - 4"                                      |
|                           |  |       |  |   |   |   | P 4 P 23 "                                |

Client:Philip Environmental Inc. Project:16766

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# Zenon Env., unmental Laboratories - Certificate of Analysis

Page 5 of 14

| ä   | Client ID:<br>Zenon ID:<br>Date Sampled:             |                      | Method<br>Blank<br>038764 96<br>96/09/23                    | Blank<br>Spike 1<br>038764 96<br>96/09/23 | Blank<br>Spike 1<br>038764 96<br>96/09/23 | Blank<br>Spike 2<br>038764 96<br>96/09/23 | Blank<br>Spike 2<br>038764 96<br>96/09/23 |
|---|--|----------------------|---|---|---|---|---|
| Component   | MDL  | Units                |   |   | % Recovery                                |   | % Recovery                                |
| Semi - volatiles via SW846  |  |                      | -0.22   |   | 4.4                                       | 7.0                                       | 70  |
| Phenol<br>Discontinue and the   | 0.11   | mg/kg                | < 0.22  | 4.4                                       | 44  | 7.0                                       | 70  |
| Bis(2-chloroethyl)ether   | 0.18   | 11                   | < 0.36  | - 4.5                                     | -<br>45                                   | 7.1                                       | - 71                                      |
| 2-Chlorophenol  | 0.27   | 14                   | <0.54   |   |   |   | 71  |
| 1,3-Dichlorobenzene   | 0.20   | '1                   | <0.40   | -   | -   | -   | -   |
| 1-4-Dichlorobenzene   | 0.20   | 11                   | <0.40   | 2.0                                       | 41  | 2.8                                       | 57  |
| 1,2-Dichlorobenzene   | 0.20   | 11                   | <0.40   | -   | -   | -   | -   |
| Bis(2-chloroisopropyl)ether   | 0.15   |                      | < 0.30  | -   | -   | -   | ~   |
| Hexachloroethane  | 0.20   | **                   | <0.40   | -   | -   | -   | -   |
| N-Nitroso-di-N-Propylamin   |  |                      | <0.42   | 2.1                                       | 43  | 3.4                                       | 68  |
| Nitrobenzene  | 0.20   |                      | <0.40   | -   | -   | -   | -   |
| Isophorone  | 0.40   | 11                   | <0.80   | -   | -   | -   | -   |
| 2-Nitrophenol   | 0.14   | n                    | <0.28   | -   | -   | -   | -   |
| 2,4-Dimethylphenol  | 0.17   | 11                   | <0.34   | -   | -   | -   | -   |
| Bis(2-chloroethoxy)methane  |  | 11                   | <0.26   | -   | -   | -   | -   |
| 2,4-Dichlorophenol  | 0.12   | 47                   | <0.24   | -   | -   | -   | -   |
| 1,2,4-Trichlorobenzene  | 0.20   | **                   | <0.40   | 2.2                                       | 45  | 3.2                                       | 64  |
| Naphthalene   | 0.03   | 11                   | <0.06   | -   | -   | -   | -   |
| Hexachlorobutadiene   | 0.20   | 11                   | <0.40   | -   | -   | -   | -   |
| 4-Chloro-3-Methylphenol   | 0.14   | 87                   | <0.28   | 4.2                                       | 42  | 7.0                                       | 70  |
| Hexachlorocyclopentadiene   | 0.20   |                      | <0.40   | -   | -   | -   | -   |
| 2,4,6-Trichlorophenol   | 0.12   | 14                   | <0.24   | -   | -   | -   | -   |
| 2-Chloronaphthalene   | 0.09   | **                   | <0.18   | -   | -   | -   | ~   |
| Acenaphthylene  | 0.04   | **                   | <0.08   | -   | -   | -   | -   |
| Dimethyl phthalate  | 0.11   | и                    | <0.22   | -   | -   | -   | -   |
| 2,6-Dinitrotoluene  | 0.06   | "                    | < 0.12  | -   | -   | -   | -   |
| Acenaphthene  | 0.07   | **                   | < 0.14  | 2.4                                       | 47  | 3.7                                       | 74  |
| 2,4-Dinitrophenol   | 0.48   | 11                   | <0.96   | -   | -   | -   | -   |
| 2,4-Dinitrotoluene  | 0.05   | 14                   | <0.10   | 1.8                                       | 37  | 3.3                                       | 66  |
| 4-Nitrophenol   | 0.14   | п                    | <0.28   | 3.5                                       | 35  | 5.7                                       | 57  |
| Fluorene  | 0.03   | "                    | < 0.06  | -   | -   | -   | -   |
| 4-Chlorophenylphenylether   | 0.09   | "                    | < 0.18  | -   | -   | -   | -   |
| Diethyl phthalate   | 0.11   |                      | < 0.22  | ~   | -   | -   | -   |
| 4,6-Dinitro-2-methylphenol  |  | **                   | < 0.30  | -   | -   | -   | -   |
| N-Nitrosodiphenylamine  | 0.19   | н                    | < 0.38  | -   | -   | -   | -   |
| 4-Bromophenylphenylether  | 0.03   | 11                   | < 0.06  | -   | -   | -   | -   |
| Hexachlorobenzene   | 0.20   |                      | <0.40   | -   | -   | -   | -   |
| Pentachlorophenol   | 0.11   | u.                   | < 0.22  | 4.1                                       | 41  | 7.3                                       | 73  |
| Phenanthrene  | 0.03   | "                    | <0.06   | -   | -   | -   | -   |
| Anthracene  | 0.02   | н                    | <0.04   | -   | -   | -   | -   |
|   |  | 14                   |   | -   | -   | -   | -   |
|   |  | ••                   |   | -   | -   | ~   | -   |
|   |  | .,                   |   | 28  | 56  | 4.0                                       | 80  |
| -   |  |                      |   |   |   |   | υU  |
|   |  | 15                   |   | -   | -   | -   | -   |
|   |  | **                   |   | •   | -   | -   | -   |
|   |  |                      |   | -   | -   | -   | -   |
| Di-n-butyl phthalate<br>Fluoranthene<br>Pyrene<br>Benzyl butyl phthalate<br>3,3-Dichlorobenzidine<br>Benzo(a)anthracene<br>Chrysene | 0.11<br>0.02<br>0.03<br>0.06<br>0.10<br>0.02<br>0.03 | 11<br>11<br>13<br>14 | <0.22<br><0.04<br><0.06<br><0.12<br><0.20<br><0.04<br><0.06 | 2.8                                       | - 56                                      | -<br><br>-<br>-<br>-                      | )   |

11

# Zenon Environmental Laboratories - Certificate of Analysis

Page 6 of 14

|                            | Client ID:<br>Zenon ID:<br>Date Sampled: |       | Method<br>Blank<br>038764 96<br>96/09/23 | Blank<br>Spike 1<br>038764 96<br>96/09/23 | Blank<br>Spike 1<br>038764 96<br>96/09/23 | Blank<br>Spike 2<br>038764 96<br>96/09/23 | Blank<br>Spike 2<br>038764 96<br>96/09/23 |
|----------------------------|--|-------|--|---|---|---|---|
| Component                  | MDL                                      | Units |  |   | % Recovery                                |   | % Recovery                                |
| Bis(2-ethylhexyl)phthalate | 0.14                                     |       | <0.28                                    | -   | -   | -   | -   |
| Di-n-octyl phthalate       | 0.11                                     | 11    | <0.22                                    | -   | -   | -   | -   |
| Benzo(b)fluoranthene       | 0.04                                     | н     | <0.08                                    | -   | -   | -   | -   |
| Benzo(k)fluoranthene       | 0.04                                     | 11    | <0.08                                    | -   | -   | -   | -   |
| Benzo(a)pyrene             | 0.05                                     | tr.   | <0.10                                    | -   | -   | -   | -   |
| Indeno(1,2,3-cd)pyrene     | 0.06                                     | и     | <0.12                                    | -   | -   | -   | -   |
| Dibenzo(a,h)anthracene     | 0.04                                     | *1    | <0.08                                    | -   | -   | -   | -   |
| Benzo(glui)perylene        | 0.04                                     | 11    | <0.08                                    | -   | -   | -   | -   |
| N-Nitrosodimethylamine     | 1.0                                      | •7    | <2.0                                     | -   | -   | -   | -   |
| Aniline                    | 0.50                                     | *1    | <1.0                                     | -   | -   | -   | -   |
| Benzyl alcohol             | 0.50                                     |       | <1.0                                     | -   | -   | -   | -   |
| Carbazole                  | 0.50                                     | .,    | <1.0                                     | -   | -   | -   | -   |
| 2-Methylphenol             | 0.50                                     | 41    | <1.0                                     | - '                                       | -   | -   | -   |
| Benzoic acid               | 0.50                                     | н     | <1.0                                     | -   | -   | -   | -   |
| 4-Chloroaniline            | 0.50                                     | **    | <1.0                                     | -   | -   | -   | -   |
| 2-Methylnaphthalene        | 0.10                                     | "     | <0.20                                    | -   | -   | -   | -   |
| 2,4,5-Trichlorophenol      | 0.10                                     | "     | <0.20                                    | -   | -   | -   | -   |
| 2-Nitroaniline             | 0.50                                     | н     | <1.0                                     | -   | -   | -   | -   |
| 3-Nitroaniline             | 0.50                                     | 11    | <1.0                                     | -   | -   | -   | -   |
| Dibenzofuran               | 0.50                                     | ч     | <1.0                                     | -   | -   | -   | -   |
| Benzidine                  | 0.50                                     | 17    | <1.0                                     | -   | -   | -   | -   |
| 4-Nitroaniline             | 0.50                                     | n     | <1.0                                     | -   | -   | -   | -   |
| Surrogate Recoveries       |  | %     |  |   |   |   |   |
| 2-Fluorophenol             |  |       | 29                                       | 42  | 42  | 52  | 52  |
| d5-Phenol                  |  |       | 55                                       | 45  | 45  | 72  | 72  |
| d5-Nitrobenzene            |  |       | 43                                       | 40  | 40  | 61  | 61  |
| 2-Fluorobiphenyl           |  |       | 65                                       | 46  | 46  | 71  | 71  |
| 2,4.6-Tribromophenol       |  |       | 66                                       | 46  | 46  | 73  | 73  |
| d14-p-Terphenyl            |  |       | 74                                       | 70  | 70  | 83  | 83  |

11



5555 North Service Road Burlington, Ontario, Canada L7L 5H7 Tel: (905) 332-8788 Fax: (905) 332-9169

# Certificate of Analysis

### **CLIENT INFORMATION**

| Attention:    | Cory Chance               |
|---------------|---------------------------|
| Client Name:  | Philip Environmental Inc. |
| Project:      | 17039                     |
| Project Desc: | WFS Septic Leach Field    |
| Address:      | 4000 Monroe Road          |
|               | Farmington, NM            |
|               | 87401                     |
| Fax Number:   | 505 326-2388              |
| Phone Number: | 505 326-2262              |

### LABORATORY INFORMATION

| Contact:        | Ada Blythe, B.Sc., C.Chem. |
|-----------------|----------------------------|
| Project:        | AN961368                   |
| Date Received:  | 96/12/13                   |
| Date Reported:  | 97/01/17                   |
| Submission No.: | 6L0416                     |
| Sample No.:     | 055065-055069              |

NOTES: ''-' = not analysed '<' = less than Method Detection Limit (MDL) 'NA' = no data available LOQ can by determined for all analytes by multiplying the appropriate MDL X 3.33 Solids data is based on dry weight except for biota analyses. Organic analyses are not corrected for extraction recovery standards except for isotope dilution methods, (i.e. CARB 429 PAH, all PCDD/F and DBD/DBF analyses)

Methods used by Zenon are based upon those found in 'Standard Methods for the Examination of Water and Wastewater', Seventeenth Edition. Other methods are based on the principles of MISA or EPA methodologies.

All work recorded herein has been done in accordance with normal professional standards using accepted testing methodologies, quality assurance and quality control procedures except where otherwise agreed to by the client and testing company in writing. Any and all use of these test results shall be limited to the actual cost of the pertinent analysis done. There is no other warranty expressed or implied. Your samples will be retained at Zenon for a period of three weeks from receipt of data or as per contract.

#### COMMENTS:

NR = 1,4-dichlorobenzene was not recovered. There were low recoveries of this compound throughout the submission. This is likely due to the volatility of this compound.

1,2,4-trichlorobenzene also had low recoveries. Again, this loss is likely due to volatility and slightly more rigorous treatment of the sample.

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Certified by.

A division of PHILIP Analytical Services Corporation

Page 1

Page 2 of 12

Zenon Environmental Laboratories - Certificate of Analysis

|                               |         |       | Method   | Blank     | Blank     |          | BHI   | BH2      | BH3       |
|-------------------------------|---------|-------|----------|-----------|-----------|----------|-------|----------|-----------|
| Cli                           | ent ID: |       | Blank    | Spike 1   | Spike 1   | BH1-3-5  | 10-12 | 15-17    | 13-15     |
| Zenon ID:                     | ton ID: |       | 05506596 | 055065 96 | 05506596  | 05506696 |       | 05506896 | 055069 96 |
| Date Sa                       | umpled: |       | 96/12/12 | 96/12/12  | 96/12/12  | 96/12/12 |       | 96/12/12 | 96/12/12  |
| Component                     | MDL     | Units |          |           | %Recovery |          |       |          |           |
| Mercury via SW846 Method 7471 | 14      |       |          |           |           |          |       |          |           |
| Mercury                       | 0.04    | mg/kg | V        | 0.99      | 66        | V        | V     | V        | V         |
| Metals via SW846 Method 6010  |         |       |          |           |           |          |       |          |           |
| Arsenic                       | 5.0     | mg/kg | <10      | 96        | 94        | <10      | <10   | <10      | <10       |
| Selenium                      | 10      | ÷     | <20      | 100       | 92        | <20      | <20   | <20      | <20       |
| Barium                        | 0.1     | =     | <0.2     | 200       | 98        | 280      | 130   | 190      | 150       |
| Cadmium                       | 0.2     | :     | <0.4     | 94        | 93        | <0.4     | <0.4  | <0.4     | <0.4      |
| Chromium                      | S       | ÷     | <10      | 200       | 66        | <10      | <10   | <10      | <10       |
| Lead                          | 10      | ÷     | <20      | 200       | 66        | <20      | <20   | <20      | <20       |
| Silver                        | 0.5     | =     | <1.0     | 96        | 96        | <1.0     | <1.0  | <1.0     | <1.0      |

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Client: Philip Environmental Inc. Project: 17039

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## Zenon Environmental Laboratories - Certificate of Analysis

|                                      | Client ID:<br>Zenon ID:<br>Date Sampled: |       | Method<br>Blank<br>055065 96<br>96/12/12 | Blank<br>Spike 1<br>055065 96<br>96/12/12 | Blank<br>Spike 1<br>055065 96<br>96/12/12 | BH1-3-5<br>055066 96<br>96/12/12 | BH1<br>10-12<br>055067 96<br>96/12/12 | BH2<br>15-17<br>055068 96<br>96/12/12 |
|--------------------------------------|--|-------|--|---|---|----------------------------------|---------------------------------------|---------------------------------------|
| Component                            | MDL                                      | Units |  |   | %Recovery                                 |                                  |                                       |                                       |
| Volatiles via SW846 Mei              | hod 8260                                 |       |  |   | £   |                                  |                                       | .,                                    |
| Acetone                              | 0.030                                    | mg/kg | '(0.17 /                                 | 0.39                                      | (160-                                     | 4 0.18 J                         | 40.15 V                               | $_{\sim 0.14}$ V                      |
| Acrolein                             | 0.010                                    | "     | <  | 0.68                                      | 140                                       | <                                | <                                     | <                                     |
| Acrylonitrile                        | 0.010                                    | "     | <  | 0.46                                      | 92  | <                                | <                                     | <                                     |
| Benzene                              | 0.005                                    | **    | <  | 0.23                                      | 94  | <                                | <                                     | <                                     |
| Bromoform                            | 0.010                                    | "     | <  | 0.26                                      | 100                                       | <                                | <                                     | <                                     |
| Bromomethane                         | 0.010                                    | ••    | <  | 0.35                                      | (140)                                     | <                                | <                                     | <                                     |
| 2-Butanone                           | 0.015                                    | **    | <  | 0.23                                      | 92  | <                                | <                                     | <                                     |
| Carbon Disulfide                     | 0.010                                    | "     | <  | 0.25                                      | 98  | <                                | <                                     | <                                     |
| Carbon Tetrachloride                 | 0.010                                    | "     | <  | 0.25                                      | 100                                       | <                                | <                                     | <                                     |
| Chlorobenzene                        | 0.005                                    | "     | <  | 0.24                                      | 97  | <                                | <                                     | <                                     |
| Chlorodibromomethane                 | 0.005                                    | ••    | <  | 0.25                                      | 100                                       | <                                | <                                     | <                                     |
| Chloroethane                         | 0.010                                    |       | <  | 0.25                                      | 100                                       | <                                | <                                     | <                                     |
| 2-Chloroethylvinylether              | 0.010                                    | "     | <  | 0.22                                      | 89  | <                                | <                                     | <                                     |
| Chloroform                           | 0.005                                    | ••    | <  | 0.22                                      | 94  | <                                | <                                     | <                                     |
| Chloromethane                        | 0.010                                    | **    | <  | 0.28                                      | 110                                       | <                                | <                                     | <                                     |
| 1,2-Dichlorobenzene                  | 0.005                                    | "     | <  | 0.25                                      | 99  | <                                | <                                     | <                                     |
| 1,3-Dichlorobenzene                  | 0.005                                    | ••    | <  | 0.25                                      | 100                                       | <                                | <                                     | <                                     |
| 1,4-Dichlorobenzene                  | 0.005                                    |       | <  | 0.26                                      | 100                                       | 0.005                            | <                                     | <                                     |
| Dichlorobromomethane                 | 0.005                                    |       | <  | 0.25                                      | 100                                       | <                                | <                                     | <                                     |
| 1,1-Dichloroethane                   | 0.005                                    |       | <  | 0.25                                      | 100                                       | <                                | <                                     | <                                     |
| ,2-Dichloroethane                    | 0.005                                    | **    | <  | 0.23                                      | 92  | <                                | <                                     | <                                     |
| ,1-Dichloroethene                    | 0.010                                    | ••    | <  | 0.25                                      | 100                                       | <                                | <                                     | <                                     |
| is-1,2-Dichloroethene                | 0.010                                    | **    | <  | 0.23                                      | 92  | <                                | <                                     | <                                     |
| rans-1,2-Dichloroethene              | 0.010                                    | 11    | <  | 0.25                                      | 99  | <                                | <                                     | <                                     |
| 1,2-Dichloropropane                  | 0.005                                    |       | <  | 0.23                                      | 91  | <                                | <                                     | <                                     |
| cis-1,3-Dichloropropene              | 0.005                                    |       | <  | 0.23                                      | 95  | <                                | <                                     | <                                     |
| rans-1,3-Dichloropropen              |  | **    | <  | 0.24                                      | 97  | <                                | <                                     | <                                     |
| Ethylbenzene                         | 0.005                                    |       | <  | 0.24                                      | 97  | <                                | <                                     | <                                     |
| 2-Hexanone                           | 0.005                                    | .,    | <  | 0.24                                      | 88  | <                                | <                                     | <                                     |
| Methylene Chloride                   | 0.010                                    | **    | 70.020                                   | 0.22                                      | 93  | < 0.039 U                        | ∠0.034 V                              | د0.029 <i>U</i>                       |
| -Methyl-2-Pentanone                  | 0.010                                    | **    | < 0.020                                  | 0.23                                      | 88  | < 0.039 C                        | <                                     | <                                     |
| Styrene                              | 0.010                                    | +1    | <  | 0.22                                      | 98  | <                                | <                                     | <                                     |
| styrene<br>1,1,1,2-Tetrachloroethane |  |       | <  | 0.24                                      | 98<br>100                                 | <                                | <                                     | <                                     |
| 1,1,2,2-Tetrachloroethane            |  |       |  | 0.23                                      | 94  | <                                | <                                     | <                                     |
| Fetrachloroethene                    | 0.010                                    | 11    | <  | 0.25                                      | 94<br>100                                 | <                                | <                                     | <                                     |
| Foluene                              | 0.020                                    | **    | <  | 0.20                                      | 96  | 0.005                            | <                                     | <                                     |
|                                      |  |       | <  | 0.24<br>0.25                              | 96<br>99                                  |                                  | <                                     | <                                     |
| 1,1,1-Trichloroethane                | 0.005<br>0.010                           | 11    | <  | 0.25                                      | 99<br>97                                  | <                                | <                                     | <                                     |
| 1,1,2-Trichloroethane                |  | ••    | <  | 0.24<br>0.24                              | 97<br>97                                  | <                                | <                                     | <                                     |
| Frichloroethene                      | 0.005                                    |       | <  |   |   | <                                |                                       | <                                     |
| Vinyl Acetate                        | 0.010                                    | **    | <  | 0.28                                      | 110                                       | <                                | <                                     |                                       |
| Vinyl Chloride                       | 0.010                                    |       | <  | 0.24                                      | 96<br>04                                  | <                                | <                                     | <                                     |
| Xylenes(Total)                       | 0.005                                    |       | <  | 0.71                                      | 94  | <                                | <                                     | <                                     |
| Surrogate Recoveries                 |  | %     | 05                                       | 0.0                                       | 00  | 07                               | 102                                   | 0.9                                   |
| d4-1,2-Dichloroethane                |  |       | 95                                       | 98  | 98  | 97                               | 103                                   | 98                                    |
| d8-Toluene                           |  |       | 98                                       | 98  | 98  | 99                               | 94<br>05                              | 98                                    |
| Bromofluorobenzene                   |  |       | 95                                       | 112                                       | 112                                       | 95                               | 95                                    | 91<br>(1)                             |

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Client:Philip Environmental Inc. Project:17039

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## Zenon Environmental Laboratories - Certificate of Analysis

|                           | Client ID:<br>Zenon ID: |       | BH3<br>13-15<br>055069 96 |
|---------------------------|-------------------------|-------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|
|                           | Date Sampled:           |       | 96/12/12                  | 96/12/12                  | 96/12/12                  | 96/12/12                  | 96/12/12                  |
| Component                 | MDL                     | Units |                           | M. Spike                  | MS % Rec.                 | MS Dup                    | MSD % Rec                 |
| Volatiles via SW846 Me    | thod 8260               |       |                           |                           | _                         |                           |                           |
| Acetone                   | 0.030                   | mg/kg | ∠ 0.12 V                  | 0.41                      | (160)                     | 0.33                      | 130                       |
| Acrolein                  | 0.010                   | **    | <                         | 0.42                      | 83                        | 0.36                      | 73                        |
| Acrylonitrile             | 0.010                   | 11    | <                         | 0.40                      | 80                        | 0.39                      | 78                        |
| Benzene                   | 0.005                   | 11    | <                         | 0.23                      | 94                        | 0.23                      | 92                        |
| Bromoform                 | 0.010                   | "     | <                         | 0.26                      | 110                       | 0.26                      | 100                       |
| Bromomethane              | 0.010                   | "     | <                         | 0.29                      | 120                       | 0.29                      | 120                       |
| 2-Butanone                | 0.015                   | 11    | <                         | 0.21                      | 83                        | 0.19                      | 75                        |
| Carbon Disulfide          | 0.010                   | 11    | <                         | 0.22                      | 88                        | 0.20                      | 82                        |
| Carbon Tetrachloride      | 0.010                   |       | <                         | 0.26                      | 110                       | 0.26                      | 100                       |
| Chlorobenzene             | 0.005                   |       | <                         | 0.23                      | 91                        | 0.23                      | 92                        |
| Chlorodibromomethane      | 0.005                   |       | <                         | 0.25                      | 100                       | 0.25                      | 100                       |
| Chloroethane              | 0.010                   | ••    | <                         | 0.27                      | 110                       | 0.25                      | 98                        |
| 2-Chloroethylvinylether   | 0.010                   | **    | <                         | 0.16                      | 64                        | 0.15                      | 60                        |
| Chloroform                | 0.005                   | u     | <                         | 0.24                      | 95                        | 0.23                      | 91                        |
| Chloromethane             | 0.010                   |       | <                         | 0.27                      | 110                       | 0.24                      | 94                        |
| 1,2-Dichlorobenzene       | 0.005                   | ••    | <                         | 0.24                      | 94                        | 0.23                      | 91                        |
| 1,3-Dichlorobenzene       | 0.005                   | н     | <                         | 0.22                      | 88                        | 0.21                      | 86                        |
| 1,4-Dichlorobenzene       | 0.005                   | ••    | <                         | 0.22                      | 87                        | 0.22                      | 88                        |
| Dichlorobromomethane      | 0.005                   | **    | <                         | 0.25                      | 100                       | 0.25                      | 100                       |
| 1,1-Dichloroethane        | 0.005                   | н     | <                         | 0.23                      | 93                        | 0.22                      | 90                        |
| 1,2-Dichloroethane        | 0.005                   | 11    | <                         | 0.24                      | 97                        | 0.23                      | 91                        |
| 1,1-Dichloroethene        | 0.010                   | 17    | <                         | 0.24                      | 97                        | 0.22                      | 89                        |
| cis-1,2-Dichloroethene    | 0.010                   |       | <                         | 0.22                      | 89                        | 0.21                      | 85                        |
| trans-1,2-Dichloroethene  | 0.010                   | 14    | <                         | 0.22                      | 88                        | 0.21                      | 84                        |
| 1,2-Dichloropropane       | 0.005                   | 15    | <                         | 0.23                      | 92                        | 0.24                      | 97                        |
| cis-1,3-Dichloropropene   | 0.005                   | 11    | <                         | 0.23                      | 90                        | 0.23                      | 92                        |
| trans-1,3-Dichloropropen  |                         | .,    | <                         | 0.22                      | 89                        | 0.23                      | 94                        |
| Ethylbenzene              | 0.005                   | **    | <                         | 0.23                      | 91                        | 0.23                      | 90                        |
| 2-Hexanone                | 0.010                   | **    | <                         | 0.18                      | 73                        | 0.18                      | 70                        |
| Methylene Chloride        | 0.010                   | *1    | ∠ 0.033 U                 | 0.25                      | 99                        | 0.24                      | 95                        |
| 4-Methyl-2-Pentanone      | 0.010                   | 0     | <                         | 0.20                      | 80                        | 0.19                      | 77                        |
| Styrene                   | 0.005                   | **    | <                         | 0.22                      | 88                        | 0.22                      | 89                        |
| 1,1,1,2-Tetrachloroethane |                         | "     | <                         | 0.26                      | 100                       | 0.26                      | 100                       |
| 1,1,2,2-Tetrachloroethane |                         | **    | <                         | 0.15                      | 61                        | 0.14                      | 57                        |
| Tetrachloroethene         | 0.020                   | н     | <                         | 0.24                      | 97                        | 0.25                      | 99                        |
| Toluene                   | 0.005                   | 11    | <                         | 0.23                      | 94                        | 0.24                      | 95                        |
| 1,1,1-Trichloroethane     | 0.005                   |       | <                         | 0.25                      | 100                       | 0.25                      | 100                       |
| 1,1,2-Trichloroethane     | 0.010                   | **    | <                         | 0.23                      | 94                        | 0.23                      | 93                        |
| Trichloroethene           | 0.005                   |       | <                         | 0.31                      | 1302                      | 0.31                      | 120                       |
| Vinýl Acetate             | 0.010                   | н     | < UT                      | 0.10                      | (A)                       | 0.072                     | 28                        |
| Vinyl Chloride            | 0.010                   | "     | <                         | 0.24                      | 95                        | 0.21                      | 84                        |
| Xylenes(Total)            | 0.005                   | ••    | <                         | 0.68                      | 91                        | 0.68                      | 90                        |
| Surrogate Recoveries      | 0.005                   | %     |                           | 0,00                      | <i>~</i> <b>1</b>         | 0,00                      | 20                        |
| d4-1,2-Dichloroethane     |                         |       | 99                        | 99                        | 99                        | 95                        | 95                        |
|                           |                         |       |                           |                           |                           |                           |                           |
| d8-Toluene                |                         |       | 95                        | 99                        | 99                        | 101                       | 101                       |

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| Component                        | Client ID:<br>Zenon ID:<br>& Sampled:<br>MDL | Units | Method<br>Blank<br>055065 96<br>96/12/12 | Blank<br>Spike 1<br>055065 96<br>96/12/12 | Blank<br>Spike 1<br>055065 96<br>96/12/12<br>%Recovery | Blank<br>Spike 2<br>055065 96<br>96/12/12 | Blank<br>Spike 2<br>055065 96<br>96/12/12<br>%Recovery |
|----------------------------------|--|-------|--|---|--|---|--|
| Semi -Volatiles via SW846        | Method 8270                                  |       |  |   |  |   |  |
| Phenol                           | 0.11   | mg/kg | < 0.22                                   | 39  | 39   | 52  | 52   |
| Bis(2-chloroethyl)ether          | 0.18   | "     | < 0.36                                   | -   | ÷  | -   | -  |
| 2-Chlorophenol                   | 0.27   | **    | < 0.54                                   | 19  | (19)   | 31  | 31   |
| 1,3-Dichlorobenzene              | 0.20   | **    | <0.40                                    | -   |  | -   | -  |
| 1-4-Dichlorobenzene              | 0.20   | "     | <0.40                                    | NR  | (NR)   | 0.54                                      | <b>A</b> 1.0   |
| 1,2-Dichlorobenzene              | 0.20   | er.   | <0.40                                    | -   | -  | -   | -  |
| Bis(2-chloroisopropyl)ether      | 0.15   | **    | < 0.30                                   | -   | -  | -   | -  |
| Hexachloroethane                 | 0.20   | "     | <0.40                                    | -   | -  | -   | -  |
| N-Nitroso-di-N-Propylamite       | . 0.21                                       | **    | <0.42                                    | 27  | 54   | 32  | 64   |
| Nitrobenzene                     | 0.20   |       | <0.40                                    | -   | -  | -   | -  |
| Isophorone                       | 0.40   | **    | < 0.80                                   | -   | -  | -   | -  |
| 2-Nitrophenol                    | 0.14   | *1    | < 0.28                                   | _   | -  | -   | -  |
| 2,4-Dimethylphenol               | 0.17   | **    | < 0.34                                   | -   | -  | -   | -  |
| Bis(2-chloroethoxy)methane       | 0.13   | **    | < 0.26                                   | -   | -  | -   | -  |
| 2,4-Dichlorophenol               | 0.12   | "     | < 0.24                                   | -   | -  | -   | -  |
| 1,2,4-Trichlorobenzene           | 0.20   | **    | < 0.40                                   | 1.8                                       | (4.0)  | 8.3                                       | $\bigcirc$   |
| Naphthalene                      | 0.03   | 0     | < 0.06                                   | -   | -  | -   | -  |
| Hexachlorobutadiene              | 0.20   | **    | <0.40                                    | -   | -  | -   | -  |
| 4-Chloro-3-Methylphenol          | 0.14   |       | < 0.28                                   | 58  | 58   | 64  | 64   |
| Hexachlorocyclopentadiene        | 0.20   | ••    | <0.40                                    | _   | -  | -   | -  |
| 2,4,6-Trichlorophenol            | 0.12   | 15    | <0.24                                    | -   | -  | -   | -  |
| 2-Chloronaphthalene              | 0.09   | *1    | < 0.18                                   | -   | -  | -   | -  |
| Acenaphthylene                   | 0.04   | **    | < 0.08                                   | -   | -  | -   | -  |
| Dimethyl phthalate               | 0.11   | 11    | < 0.22                                   | -   | -  | -   | -  |
| 2,6-Dinitrotoluene               | 0.06   | "     | < 0.12                                   | -   | -  | -   | -  |
| Acenaphthene                     | 0.07   | "     | < 0.14                                   | 29  | 58   | 31  | 62   |
| 2,4-Dinitrophenol                | 0.48   | **    | <0.96                                    | -   | _  | _   | _  |
| 2,4-Dinitrotoluene               | 0.05   | U.    | < 0.10                                   | 32  | 63   | 31  | 62   |
| 4-Nitrophenol                    | 0.14   | "     | <0.28                                    | 32<br>75                                  | 75   | 68  | 68   |
| Fluorene                         | 0.03   | 11    | <0.06                                    | -   | -  | -   | -  |
| 4-Chlorophenylphenylether        | 0.09   | u.    | <0.18                                    | _   | _  | -   | -  |
| Diethyl phthalate                | 0.11   |       | <0.22                                    | _   | _  | _   | -  |
| 4,6-Dinitro-2-methylphenol       | 0.15   | 11    | < 0.30                                   | _   | _  | _   | -  |
| N-Nitrosodiphenylamine           | 0.15   | "     | <0.38                                    | -   | -  | -   | _  |
| 4-Bromophenylphenylether         | 0.03   | п     | <0.08                                    | -   | -  | -   | -  |
| Hexachlorobenzene                | 0.03   | "     | <0.00                                    | ~   | -  | -   | -  |
| Pentachlorophenol                | 0.20   |       | <0.40                                    | - 69                                      | - 69   | 64  | 64   |
| Phenanthrene                     | 0.11   |       | <0.22                                    | 09  | 09   | 04  |  |
| Anthracene                       |  | "     |  | -   | -  | -   | -  |
|                                  | 0.02   | н     | <0.04                                    | -   | -  | -   | -  |
| Di-n-butyl phthalate             | 0.11   | **    | <0.22                                    | -   | -  | -   | -  |
| Fluoranthene                     | 0.02   | **    | <0.04                                    | -   | -<br>77  | -   | -  |
| Pyrene<br>Bangud butul phthalatu | 0.03   | *1    | <0.06                                    | 36  | 73   | 34  | 68   |
| Benzyl butyl phthalate           | 0.06   |       | < 0.12                                   | -   | -  | -   | -  |

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| Zenon ID:<br>Zenon ID:<br>Date Sampled:055065 96 <th< th=""><th></th><th>Client ID:</th><th></th><th>Method<br/>Blank</th><th>Blank<br/>Spike 1</th><th>Blank<br/>Spike 1</th><th>Blank<br/>Spike 2</th><th>Blank<br/>Spike 2</th></th<> |  | Client ID:  |  | Method<br>Blank  | Blank<br>Spike 1                   | Blank<br>Spike 1      | Blank<br>Spike 2                   | Blank<br>Spike 2 |   |   |   |   |
|---|--|---|--|--|------------------------------------|-----------------------|------------------------------------|------------------|---|---|---|---|
| Date Sampled:96/12/1296/1296/1296/1296/12   |  |   |  |  | -                                  | •                     | •                                  |                  |   |   |   |   |
| Component         MDL         Units $\%$ Recovery         %Recovery         %Recovery           3.3-Dichlorobenzidine         0.10 $< 0.02$ $  -$ Benzo(a)anthracene         0.02         "< $<0.04$ $  -$ Benzo(a)anthracene         0.03         "< $<0.06$ $  -$ Bis(2-ethythexyl)phthalate         0.14         "< $<0.28$ $  -$ Benzo(k)fluoranthene         0.04         "< $<0.28$ $  -$ Benzo(k)fluoranthene         0.04         "< $<0.08$ $  -$ Benzo(a)pyrene         0.06         "< $<0.12$ $  -$ Dibenzo(a, h)anthracene         0.04         "< $<0.08$ $  -$ Benzo(a)pyrene         0.06         "< $<0.08$ $  -$ N-Nitrosodimethylanine         1.0         "< $<2.0$ $  -$ Benzo(a)piorylene         0.50         "< $<1.0$ $  -$  | D  |   |  |  |                                    |                       |                                    |                  |   |   |   |   |
| 3.3-Dichlorobenzidine       0.10 $< 0.20$ $   -$ Benzo(a)anthracene       0.02 $< 0.04$ $   -$ Chrysene       0.03 $< 0.06$ $   -$ Di-n-octyl phthalate       0.11 $< 0.22$ $  -$ Benzo(b)fluoranthene       0.04 $< 0.08$ $  -$ Benzo(a)pyrene       0.05 $< 0.010$ $  -$ Benzo(a)pyrene       0.06 $< 0.12$ $  -$ Dibenzo(a,h)anthracene       0.04 $< 0.08$ $  -$ Dibenzo(a,h)anthracene       0.04 $< 0.08$ $  -$ Dibenzo(a,h)anthracene       0.04 $< < 2.00$ $  -$ N-Nitrosodimethylamine       1.0 $< < 2.0$ $  -$ Benzoic acid       0.50 $< < 1.0$ $  -$ Benzoic acid       0.50 $< < 1.0$ $ -$  |  | •   | Units  | >0/12/12   | <i>J</i> ( <i>j</i> 12 <i>j</i> 12 |                       | <i>y</i> 0 <i>/</i> 1 <i>2/</i> 12 |                  |   |   |   |   |
| Benzo(a)anthracene $0.02$ $<0.04$ $  -$ Chrysene $0.03$ $<0.06$ $  -$ Bis(2-ethylhexyl)phthalate $0.14$ $<0.28$ $  -$ Di-n-octyl phthalate $0.11$ $<0.22$ $  -$ Benzo(k)fluoranthene $0.04$ $<0.08$ $  -$ Benzo(a)pyrene $0.04$ $<0.08$ $  -$ Benzo(a)pyrene $0.06$ $<0.12$ $  -$ Benzo(a)pyrene $0.06$ $<0.08$ $   -$ Dibenzo(a,h)authracene $0.04$ $<0.08$ $   -$ Benzo(a)hjorylene $0.04$ $<0.08$ $   -$ Dibenzo(a,h)authracene $0.04$ $<0.08$ $   -$ Benzo(a)hjorylene $0.04$ $<0.08$ $    -$ Varitrosodimethylamine $1.0$ <  | -  |   |  | <0.20  | _                                  | -                     | -                                  | -                |   |   |   |   |
| $\begin{array}{cccccccccccccccccccccccccccccccccccc$  |  |   | **   |  |                                    | -                     | _                                  | -                |   |   |   |   |
| Bis(2-ethylhexyl)phthalate       0.14       " $< 0.28$ -       -       -       -         Di-n-octyl phthalate       0.11       " $< 0.22$ -       -       -       -         Benzo(k)fluoranthene       0.04       " $< 0.08$ -       -       -       -         Benzo(k)fluoranthene       0.04       " $< 0.08$ -       -       -       -         Benzo(a)pyrene       0.05       " $< 0.10$ -       -       -       -       -         Indeno(1,2,3-cd)pyrene       0.06       " $< 0.12$ -         |  |   | **   |  | -                                  | _                     | _                                  | -                |   |   |   |   |
| Di-n-octyl phthalate       0.11       "       <0.22       -       -       -       -         Benzo(b)fluoranthene       0.04       "       <0.08   | -  |   | "  |  | -                                  | -                     | -                                  | -                |   |   |   |   |
| Benzo(b)Ituoranthene $0.04$ " $<0.08$ $  -$ Benzo(a)pyrene $0.05$ " $<0.10$ $  -$ Indeno(1,2,3-cd)pyrene $0.06$ " $<0.12$ $  -$ Dibenzo(a,h)anthracene $0.04$ " $<0.08$ $  -$ Benzo(ghi)perylene $0.04$ " $<0.08$ $  -$ N-Nitrosodimethylamine $1.0$ " $<2.0$ $  -$ Aniline $0.50$ " $<1.0$ $   -$ Benzoighi)perylenol $0.50$ " $<1.0$ $   -$ N-Nitrosodimethylamine $0.50$ "<<1.0  |  |   | н  |  | -                                  | -                     | _                                  | _                |   |   |   |   |
| Benzo(k)fluoranthene $0.04$ " $< 0.08$ -       -       -       -         Benzo(a)pyrene $0.05$ " $< 0.10$ -       -       -       -         Indeno(1,2,3-cd)pyrene $0.06$ " $< 0.12$ -       -       -       -         Dibenzo(a,h)anthracene $0.04$ " $< 0.08$ -       -       -       -         Benzo(ghi)perylene $0.04$ " $< 0.08$ -       -       -       -         N-Nitrosodimethylamine $1.0$ " $< 2.0$ -       -       -       -         Aniline $0.50$ " $< 1.0$ -       -       -       -         Benzola cid $0.50$ " $< 1.0$ -       -       -       -         Carbazole $0.50$ " $< 1.0$ -       -       -       -         2-Methylphenol $0.50$ " $< 1.0$ -       -       -       -         2-Methylphenol $0.50$ "<   |  |   | 11   |  | -                                  | -                     | -                                  | -                |   |   |   |   |
| Benzo(a)pyrene $0.05$ " $<0.10$ -         |  |   | n  |  | -                                  | -                     | -                                  | -                |   |   |   |   |
| Indeno(1,2,3-cd)pyrene0.06"<0.12Dibenzo(a,h)anthracene0.04"<0.08  |  | 0.05  | "  |  | -                                  | -                     | -                                  | -                |   |   |   |   |
| Dibenzo(a,h)anthracene $0.04$ " $<0.08$ Benzo(ghi)perylene $0.04$ " $<0.08$ N-Nitrosodimethylamine $1.0$ " $<2.0$ Aniline $0.50$ " $<1.0$ Benzyl alcohol $0.50$ " $<1.0$ Carbazole $0.50$ " $<1.0$ 2-Methylphenol $0.50$ " $<1.0$ Benzoic acid $0.50$ " $<1.0$ 4-Chloroaniline $0.50$ " $<1.0$ 2-Methylnaphthalene $0.10$ "<  |  | 0.06  | 11   |  | -                                  | -                     | -                                  | -                |   |   |   |   |
| DelizionDelizionDelizionColorColo   |  | 0.04  | **   | < 0.08   | -                                  | -                     | -                                  | -                |   |   |   |   |
| A-Nitrosoumentry annue1.0 $< 2.0$ $    -$ Aniline $0.50$ " $<1.0$ $    -$ Benzyl alcohol $0.50$ " $<1.0$ $   -$ Carbazole $0.50$ " $<1.0$ $   -$ 2-Methylphenol $0.50$ " $<1.0$ $   -$ Benzoic acid $0.50$ " $<1.0$ $   -$ 4-Chloroaniline $0.50$ " $<1.0$ $  -$ 2-Methylnaphthalene $0.10$ " $<0.20$ $  -$ 2-Methylnaphthalene $0.10$ " $<0.20$ $  -$ 2-Mitroaniline $0.50$ "< <td><math>&lt;1.0</math><math>  -</math>2-Nitroaniline<math>0.50</math>"&lt;<td><math>&lt;1.0</math><math>  -</math>3-Nitroaniline<math>0.50</math>"&lt;<td><math>&lt;1.0</math><math>  -</math>Dibenzofuran<math>0.50</math>"&lt;<td><math>&lt;1.0</math><math>  -</math>Surrogate Recoveries<math>\%</math><math>55</math><math>11</math><math>11</math><math>22</math><math>22</math>d5-Phenol<math>58</math><math>31</math><math>31</math><math>51</math><math>51</math>d5-Nitrobenzene<math>\%</math><math>73</math><math>41</math><math>41</math><math>59</math><math>59</math><math>2.4.6-Tribromophenol<math>101</math><math>64</math><math>64</math><math>65</math><math>65</math></math></td><td></td><td>0.04</td><td>11</td><td>&lt; 0.08</td><td>-</td><td>-</td><td>-</td><td>-</td></td></td></td>   | $<1.0$ $  -$ 2-Nitroaniline $0.50$ "< <td><math>&lt;1.0</math><math>  -</math>3-Nitroaniline<math>0.50</math>"&lt;<td><math>&lt;1.0</math><math>  -</math>Dibenzofuran<math>0.50</math>"&lt;<td><math>&lt;1.0</math><math>  -</math>Surrogate Recoveries<math>\%</math><math>55</math><math>11</math><math>11</math><math>22</math><math>22</math>d5-Phenol<math>58</math><math>31</math><math>31</math><math>51</math><math>51</math>d5-Nitrobenzene<math>\%</math><math>73</math><math>41</math><math>41</math><math>59</math><math>59</math><math>2.4.6-Tribromophenol<math>101</math><math>64</math><math>64</math><math>65</math><math>65</math></math></td><td></td><td>0.04</td><td>11</td><td>&lt; 0.08</td><td>-</td><td>-</td><td>-</td><td>-</td></td></td> | $<1.0$ $  -$ 3-Nitroaniline $0.50$ "< <td><math>&lt;1.0</math><math>  -</math>Dibenzofuran<math>0.50</math>"&lt;<td><math>&lt;1.0</math><math>  -</math>Surrogate Recoveries<math>\%</math><math>55</math><math>11</math><math>11</math><math>22</math><math>22</math>d5-Phenol<math>58</math><math>31</math><math>31</math><math>51</math><math>51</math>d5-Nitrobenzene<math>\%</math><math>73</math><math>41</math><math>41</math><math>59</math><math>59</math><math>2.4.6-Tribromophenol<math>101</math><math>64</math><math>64</math><math>65</math><math>65</math></math></td><td></td><td>0.04</td><td>11</td><td>&lt; 0.08</td><td>-</td><td>-</td><td>-</td><td>-</td></td> | $<1.0$ $  -$ Dibenzofuran $0.50$ "< <td><math>&lt;1.0</math><math>  -</math>Surrogate Recoveries<math>\%</math><math>55</math><math>11</math><math>11</math><math>22</math><math>22</math>d5-Phenol<math>58</math><math>31</math><math>31</math><math>51</math><math>51</math>d5-Nitrobenzene<math>\%</math><math>73</math><math>41</math><math>41</math><math>59</math><math>59</math><math>2.4.6-Tribromophenol<math>101</math><math>64</math><math>64</math><math>65</math><math>65</math></math></td> <td></td> <td>0.04</td> <td>11</td> <td>&lt; 0.08</td> <td>-</td> <td>-</td> <td>-</td> <td>-</td> | $<1.0$ $  -$ Surrogate Recoveries $\%$ $55$ $11$ $11$ $22$ $22$ d5-Phenol $58$ $31$ $31$ $51$ $51$ d5-Nitrobenzene $\%$ $73$ $41$ $41$ $59$ $59$ $2.4.6-Tribromophenol10164646565$ |                                    | 0.04                  | 11                                 | < 0.08           | - | - | - | - |
| Benzyl alcohol $0.50$ " $<1.0$ Carbazole $0.50$ " $<1.0$ 2-Methylphenol $0.50$ " $<1.0$ Benzoic acid $0.50$ " $<1.0$ 4-Chloroaniline $0.50$ " $<1.0$ 2-Methylnaphthalene $0.10$ " $<0.20$ 2.4,5-Trichlorophenol $0.10$ " $<0.20$ 2.4,5-Trichlorophenol $0.50$ " $<1.0$ 2.Nitroaniline $0.50$ " $<1.0$ Dibenzofuran $0.50$ " $<1.0$ Benzidine $0.50$ " $<1.0$ Surrogate Recoveries% $<$  | N-Nitrosodimethylamine   | 1.0   |  | <2.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| Carbazole $0.50$ " $<1.0$ 2-Methylphenol $0.50$ " $<1.0$ Benzoic acid $0.50$ " $<1.0$ 4-Chloroaniline $0.50$ " $<1.0$ 2-Methylnaphthalene $0.10$ " $<0.20$ 2-Methylnaphthalene $0.10$ " $<0.20$ 2.4,5-Trichlorophenol $0.10$ " $<0.20$ 2-Nitroaniline $0.50$ " $<1.0$ 3-Nitroaniline $0.50$ "<  | Aniline  | 0.50  | 11   | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| $\begin{array}{cccccccccccccccccccccccccccccccccccc$  | Benzyl alcohol   | 0.50  | н  | <1.0   | -                                  | ~                     | -                                  | -                |   |   |   |   |
| Benzoic acid $0.50$ " $<1.0$ 4-Chloroaniline $0.50$ " $<1.0$ 2-Methylnaphthalene $0.10$ " $<0.20$ 2.4,5-Trichlorophenol $0.10$ " $<0.20$ 2-Nitroaniline $0.50$ " $<1.0$ 3-Nitroaniline $0.50$ " $<1.0$ Dibenzofuran $0.50$ " $<1.0$ Benzidine $0.50$ " $<1.0$ 4-Nitroaniline $0.50$ "< <td><math>&lt;1.0</math>Surrogate Recoveries%<math><math><math><math><math><math><math><math>2-Fluorophenol<math><math><math><math>&lt;</math></math></math></math></math></math></math></math></math></math></math></math></td> $<$   | $<1.0$ Surrogate Recoveries% $2-Fluorophenol<$   | Carbazole   | 0.50   | 17   | <1.0                               | -                     | -                                  | -                | - |   |   |   |
| 4-Chloroaniline $0.50$ " $<1.0$ 2-Methylnaphthalene $0.10$ " $<0.20$ $2,4,5$ -Trichlorophenol $0.10$ " $<0.20$ 2-Nitroaniline $0.50$ " $<1.0$ 3-Nitroaniline $0.50$ " $<1.0$ Dibenzofuran $0.50$ " $<1.0$ Benzidine $0.50$ " $<1.0$ 4-Nitroaniline $0.50$ " $<1.0$ Surrogate Recoveries% $<25$ $11$ $(11)$ $22$ $(22)$ d5-Phenol $58$ $31$ $31$ $51$ $51$ d5-Nitrobenzene $\sqrt{73}$ $41$ $41$ $59$ $59$ 2.4,6-Tribromophenol $101$ $64$ $64$ $65$ $65$  | 2-Methylphenol   | 0.50  | 18   | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| $\begin{array}{cccccccccccccccccccccccccccccccccccc$  | Benzoic acid   | 0.50  | "  | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| 2,4,5-Trichlorophenol0,10"<0.202-Nitroaniline0,50"<1.0  | 4-Chloroaniline  | 0.50  | 11   | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| 2,4,5) Themiolophend0.10 $\langle 0.20$ $i$ $i$ $i$ $i$ $i$ 2-Nitroaniline $0.50$ " $\langle 1.0$ $   -$ 3-Nitroaniline $0.50$ " $\langle 1.0$ $   -$ Dibenzofuran $0.50$ " $\langle 1.0$ $   -$ Benzidine $0.50$ " $\langle 1.0$ $   -$ 4-Nitroaniline $0.50$ " $\langle 1.0$ $  -$ Surrogate Recoveries% $\langle 25 \rangle$ $11$ $11$ $22$ $22$ d5-Phenol $58$ $31$ $31$ $51$ $51$ d5-Nitrobenzene $\gamma$ -tv $(19)$ $12$ $(12)$ $30$ $30$ 2-Fluorobiphenyl $73$ $41$ $41$ $59$ $59$ 2.4,6-Tribromophenol $101$ $64$ $64$ $65$ $65$   | 2-Methylnaphthalene  | 0.10  | "  | <0.20  | -                                  | -                     | -                                  | -                |   |   |   |   |
| 2-Mitoannic $0.50$ $<1.0$ $   -$ 3-Nitroaniline $0.50$ " $<1.0$ $   -$ Dibenzofuran $0.50$ " $<1.0$ $   -$ Benzidine $0.50$ " $<1.0$ $   -$ 4-Nitroaniline $0.50$ " $<1.0$ $   -$ Surrogate Recoveries $\%$ $55$ $11$ $11$ $22$ $22$ 2-Fluorophenol $58$ $31$ $31$ $-51$ $51$ d5-Phenol $58$ $31$ $31$ $-51$ $51$ d5-Nitrobenzene $\rightarrow$ $19$ $12$ $12$ $12$ $30$ 2-Fluorobiphenyl $73$ $41$ $41$ $59$ $59$ 2.4,6-Tribromophenol $101$ $64$ $64$ $65$ $65$   | 2,4,5-Trichlorophenol  | 0.10  | н  | < 0.20   | -                                  | -                     | -                                  | -                |   |   |   |   |
| 3-Nitroamme $0.50$ $<1.0$ $   -$ Dibenzofuran $0.50$ " $<1.0$ $   -$ Benzidine $0.50$ " $<1.0$ $   -$ 4-Nitroaniline $0.50$ " $<1.0$ $   -$ Surrogate Recoveries% $  -2-Fluorophenol5831131<5151d5-Phenol5831131<5151d5-Nitrobenzene >+\sqrt{3}12122-Fluorobiphenyl73414159592.4,6-Tribromophenol10164646565$   | 2-Nitroaniline   | 0.50  | 11   | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| Benzidine $0.50$ " $<1.0$ 4-Nitroaniline $0.50$ " $<1.0$ Surrogate Recoveries% $\sim$ $\sim$ 2-Fluorophenol $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ d5-Phenol $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ d5-Nitrobenzene $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ d5-Nitrobenzene $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ d5-Nitrobenzene $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ $\sim$ 2-Fluorobiphenyl73414159592.4,6-Tribromophenol10164646565   | 3-Nitroaniline   | 0.50  | **   | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| 4-Nitroaniline $0.50$ "<1.0Surrogate Recoveries% $55$ 111122222-Fluorophenol $55$ 11112222d5-Phenol $58$ 31315151d5-Nitrobenzene<br>2-Fluorobiphenyl $73$ 414159592.4,6-Tribromophenol10164646565   |  | 0.50  | 11   | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| Surrogate Recoveries%111122222-Fluorophenol $58$ $31$ $31$ $51$ $51$ d5-Phenol $58$ $31$ $31$ $51$ $51$ d5-Nitrobenzene $40$ $19$ $12$ $12$ $30$ $30$ 2-Fluorobiphenyl $73$ $41$ $41$ $59$ $59$ 2.4,6-Tribromophenol $101$ $64$ $64$ $65$ $65$  | Benzidine  |   |  | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| 2-Fluorophenol $55$ 11 $11$ 22 $22$ d5-Phenol583131-5151d5-Nitrobenzene $19$ 12 $12$ $30$ 302-Fluorobiphenyl73414159592,4,6-Tribromophenol10164646565   |  | 0.50  | **   | <1.0   | -                                  | -                     | -                                  | -                |   |   |   |   |
| d5-Phenol $58$ $31$ $31$ $-51$ $51$ d5-Nitrobenzene $19$ $12$ $12$ $30$ $30$ 2-Fluorobiphenyl $73$ $41$ $41$ $59$ $59$ 2.4.6-Tribromophenol $101$ $64$ $64$ $65$ $65$   |  |   | %  | $\bigcirc$   |                                    | $\cap$                |                                    | $\frown$         |   |   |   |   |
| d5-Nitrobenzene12121230302-Fluorobiphenyl73414159592.4.6-Tribromophenol10164646565  |  |   |  | 25   |                                    | $\forall \mathcal{V}$ |                                    | (22)             |   |   |   |   |
| 2-Fluorobiphenyl73414159592,4,6-Tribromophenol10164646565   |  |   |  | 58   |                                    | 31                    |                                    |                  |   |   |   |   |
| 2,4,6-Tribromophenol 101 64 64 65 65  |  |   |  | ~  |                                    |                       |                                    |                  |   |   |   |   |
| •   | , -  |   |  |  |                                    |                       |                                    |                  |   |   |   |   |
| d14-p-Terphenyl 105 77 77 73 73   | -  |   |  |  |                                    |                       |                                    |                  |   |   |   |   |
|   | d14-p-Terphenyl  |   |  | 105  | 77                                 | 77                    | 73                                 | 73               |   |   |   |   |

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| Dat                         | Client ID:<br>Zenon ID:<br>e Sampled: |            | BH1-3-5<br>055066 96<br>96/12/12 |
|-----------------------------|---------------------------------------|------------|----------------------------------|----------------------------------|----------------------------------|----------------------------------|----------------------------------|
| Component                   | MDL                                   | Units      | ,                                | M. Spike                         | MS % Rec.                        | MS Dup                           | MSD % Rec.                       |
| Semi -Volatiles via SW846 M |                                       |            |                                  | M. Spike                         | W15 /// KCC.                     | MO Dup                           | 1013D 70 Rec.                    |
| Phenol                      | 0.11                                  | ,<br>mg/kg | < 0.22                           | 6.0                              | 67                               | 6.1                              | 59                               |
| Bis(2-chloroethyl)ether     | 0.11                                  | mg/kg      | <0.22                            | 0.0                              | 07                               | 0.1                              | .)7                              |
| 2-Chlorophenol              | 0.18                                  | "          | <0.54 <i>V</i> J                 | -<br>-<br>4.4                    | -<br>49                          | 4.9                              | 47                               |
| 1,3-Dichlorobenzene         | 0.27                                  | **         | <0.40                            | 4.4                              | 49                               |                                  | 47                               |
| 1-4-Dichlorobenzene         | 0.20                                  | н          | <0.40 R                          | 0.10                             | (2.9)                            | 0.14                             | ( <del>5</del> .0)               |
| 1,2-Dichlorobenzene         | 0.20                                  | ••         | <0.40                            | 0.10                             | 2.9                              | 0.14                             |                                  |
| Bis(2-chloroisopropyl)ether | 0.15                                  |            | <0.30                            | _                                | _                                | _                                | -                                |
| Hexachloroethane            | 0.20                                  | **         | <0.40                            | -                                | _                                | -                                | -                                |
| N-Nitroso-di-N-Propylamine  |                                       | 11         | <0.42                            | 2.8                              | 62                               | 2.8                              | 55                               |
| Nitrobenzene                | 0.20                                  |            | <0.40                            | -                                | -                                | -                                |                                  |
| Isophorone                  | 0.40                                  | "          | <0.80                            | _                                | _                                | _                                | -                                |
| 2-Nitrophenol               | 0.14                                  | "          | <0.28                            | -                                | _                                | -                                | -                                |
| 2,4-Dimethylphenol          | 0.17                                  | "          | < 0.34                           | _                                | _                                | -                                | _                                |
| Bis(2-chloroethoxy)methane  | 0.13                                  |            | <0.26                            | -                                | _                                | _                                | -                                |
| 2,4-Dichlorophenol          | 0.12                                  | "          | <0.24                            | -                                |                                  | _                                | _                                |
| 1,2,4-Trichlorobenzene      | 0.20                                  | "          | < 0.40 R                         | 0.94                             | (21)                             | 1.0                              | (19)                             |
| Naphthalene                 | 0.03                                  | 11         | <0.06                            | -                                | e -                              | -                                | <u>e</u>                         |
| Hexachlorobutadiene         | 0.20                                  | **         | <0.40                            | -                                | _                                | -                                | -                                |
| 4-Chloro-3-Methylphenol     | 0.14                                  |            | <0.28                            | 5.2                              | 58                               | 4.9                              | 47                               |
| Hexachlorocyclopentadiene   | 0.20                                  | 11         | < 0.40                           | -                                | -                                | -                                | -                                |
| 2,4,6-Trichlorophenol       | 0.12                                  |            | < 0.24                           | -                                | -                                | _                                | -                                |
| 2-Chloronaphthalene         | 0.09                                  |            | < 0.18                           | -                                | -                                | -                                | -                                |
| Acenaphthylene              | 0.04                                  | "          | < 0.08                           | -                                | -                                | _                                | -                                |
| Dimethyl phthalate          | 0.11                                  | **         | < 0.22                           | -                                | _                                | -                                | -                                |
| 2,6-Dinitrotoluene          | 0.06                                  | 11         | < 0.12                           | -                                | -                                | -                                | -                                |
| Acenaphthene                | 0.07                                  | **         | < 0.14                           | 3.1                              | 68                               | 3.1                              | 60                               |
| 2,4-Dinitrophenol           | 0.48                                  | 17         | < 0.96                           | -                                | -                                | -                                | -                                |
| 2,4-Dinitrotoluene          | 0.05                                  | "          | < 0.10                           | 2.8                              | 62                               | 3.1                              | 59                               |
| 4-Nitrophenol               | 0.14                                  | 11         | < 0.28                           | 6.0                              | 66                               | 7.1                              | 68                               |
| Fluorene                    | 0.03                                  | 11         | < 0.06                           | -                                | -                                | · _                              | -                                |
| 4-Chlorophenylphenylether   | 0.09                                  | "          | < 0.18                           | _                                | -                                | -                                | -                                |
| Diethyl phthalate           | 0.11                                  | "          | < 0.22                           | -                                | -                                | -                                | -                                |
| 4,6-Dinitro-2-methylphenol  | 0.15                                  | "          | < 0.30                           | -                                | -                                | -                                | -                                |
| N-Nitrosodiphenylamine      | 0.19                                  | "          | < 0.38                           | -                                | -                                | -                                | -                                |
| 4-Bromophenylphenylether    | 0.03                                  | н          | < 0.06                           | -                                | -                                | -                                | -                                |
| Hexachlorobenzene           | 0.20                                  | "          | < 0.40                           | -                                | -                                | -                                | -                                |
| Pentachlorophenol           | 0.11                                  | "          | < 0.22                           | 6.2                              | 69                               | 6.0                              | 78                               |
| Phenanthrene                | 0.03                                  | "          | < 0.06                           | -                                | -                                | -                                | -                                |
| Anthracene                  | 0.02                                  | "          | < 0.04                           | -                                | -                                | -                                | -                                |
| Di-n-butyl phthalate        | 0.11                                  | **         | < 0.22                           | -                                | -                                | -                                | -                                |
| Fluoranthene                | 0.02                                  | н          | < 0.04                           | -                                | -                                | -                                | -                                |
| Pyrene                      | 0.03                                  | **         | < 0.06                           | 3.3                              | 72                               | 3.5                              | 68                               |
| Benzyl butyl phthalate      | 0.06                                  | "          | < 0.12                           | -                                | -                                | -                                | -                                |

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| Da                         | Client ID:<br>Zenon ID:<br>ate Sampled: |       | BH1-3-5<br>055066 96<br>96/12/12 |
|----------------------------|---|-------|----------------------------------|----------------------------------|----------------------------------|----------------------------------|----------------------------------|
| Component                  | MDL                                     | Units |                                  | M. Spike                         | MS % Rec.                        | MS Dup                           | MSD % Rec.                       |
| 3,3-Dichlorobenzidine      | 0.10                                    | "     | <0.20                            | -                                | -                                | -                                | -                                |
| Benzo(a)anthracene         | 0.02                                    | 11    | < 0.04                           | -                                | -                                | -                                | -                                |
| Chrysene                   | 0.03                                    | "     | < 0.06                           | -                                | -                                | -                                | -                                |
| Bis(2-ethylhexyl)phthalate | 0.14                                    | н     | < 0.28                           | -                                | -                                | -                                | -                                |
| Di-n-octyl phthalate       | 0.11                                    | "     | < 0.22                           | -                                | -                                | -                                | -                                |
| Benzo(b)fluoranthene       | 0.04                                    | "     | < 0.08                           | -                                | -                                | -                                | -                                |
| Benzo(k)fluoranthene       | 0.04                                    |       | < 0.08                           | -                                | -                                | -                                | -                                |
| Benzo(a)pyrene             | 0.05                                    | "     | < 0.10                           | -                                | -                                | -                                | -                                |
| Indeno(1,2,3-cd)pyrene     | 0.06                                    | н     | < 0.12                           | -                                | -                                | -                                | -                                |
| Dibenzo(a,h)anthracene     | 0.04                                    | "     | < 0.08                           | -                                | -                                | -                                | -                                |
| Benzo(ghi)perylene         | 0.04                                    | **    | < 0.08                           | -                                | -                                | -                                | -                                |
| N-Nitrosodimethylamine     | 1.0                                     |       | <2.0                             | -                                | -                                | -                                | -                                |
| Aniline                    | 0.50                                    | **    | <1.0                             | -                                | -                                | -                                | -                                |
| Benzyl alcohol             | 0.50                                    |       | <1.0                             | -                                | -                                | -                                | -                                |
| Carbazole                  | 0.50                                    | **    | <1.0                             | -                                | -                                | -                                | -                                |
| 2-Methylphenol             | 0.50                                    | н     | <1.0                             | -                                | -                                | -                                | -                                |
| Benzoic acid               | 0.50                                    | 11    | <1.0                             | -                                | -                                | -                                | -                                |
| 4-Chloroaniline            | 0.50                                    | 18    | <1.0                             | -                                | -                                | -                                | -                                |
| 2-MethyInaphthalene        | 0.10                                    | 0     | < 0.20                           | -                                | -                                | -                                | -                                |
| 2,4,5-Trichlorophenol      | 0.10                                    | "     | < 0.20                           | -                                | -                                | -                                | -                                |
| 2-Nitroaniline             | 0.50                                    | **    | <1.0                             | -                                | -                                | -                                | -                                |
| 3-Nitroaniline             | 0.50                                    | 11    | <1.0                             | -                                | -                                | -                                | -                                |
| Dibenzofuran               | 0.50                                    | ŧſ    | <1.0                             | -                                | -                                | -                                | -                                |
| Benzidine                  | 0.50                                    | *1    | <1.0                             | -                                | -                                | -                                | -                                |
| 4-Nitroaniline             | 0.50                                    | **    | <1.0                             | -                                | -                                | -                                | -                                |
| Surrogate Recoveries       |   | %     |                                  |                                  |                                  |                                  |                                  |
| 2-Fluorophenol             |   |       | 40                               | 44                               | 44                               | 45                               | 45                               |
| d5-Phenol                  |   |       | 68                               | 69                               | 69                               | 59                               | 59                               |
| d5-Nitrobenzene            |   |       | 42                               | 46                               | 46                               | 43                               | 53                               |
| 2-Fluorobiphenyl           |   |       | 59                               | 64                               | 64                               | 54                               | 54                               |
| 2,4,6-Tribromophenol       |   |       | 69                               | 70                               | 70                               | 68                               | 68                               |
| d14-p-Terphenyl            |   |       | 75                               | 77                               | 77                               | 74                               | 74                               |

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| Z                            | Client ID:<br>enon ID:<br>Sampled:<br>MDL | Units | BH1<br>10-12<br>055067 96<br>96/12/12 | BH2<br>15-17<br>055068 96<br>96/12/12 | BH3<br>13-15<br>055069 96<br>96/12/12 |
|------------------------------|---|-------|---------------------------------------|---------------------------------------|---------------------------------------|
| Semi -Volatiles via SW846 Me | thod 8270                                 | 1     |                                       |                                       |                                       |
| Phenol                       | 0.11                                      | mg/kg | < 0.22                                | <0.22                                 | < 0.22                                |
| Bis(2-chloroethyl)ether      | 0.18                                      | "     | < 0.36                                | < 0.36                                | < 0.36                                |
| 2-Chlorophenol               | 0.27                                      | **    | <0.54 UT                              | - <0.540J                             | حرم 0.54<                             |
| 1,3-Dichlorobenzene          | 0.20                                      | "     | <0.40                                 | <0.40                                 | <0.40                                 |
| 1-4-Dichlorobenzene          | 0.20                                      |       | <0:40 R                               | <0.40L                                | 50.40 R                               |
| 1,2-Dichlorobenzene          | 0.20                                      | "     | <0.40                                 | <0.40                                 | <0.40                                 |
| Bis(2-chloroisopropyl)ether  | 0.15                                      | "     | < 0.30                                | <0.30                                 | < 0.30                                |
| Hexachloroethane             | 0.20                                      | **    | <0.40                                 | <0.40                                 | <0.40                                 |
| N-Nitroso-di-N-Propylamine   | 0.21                                      | **    | < 0.42                                | <0.42                                 | < 0.42                                |
| Nitrobenzene                 | 0.20                                      | н     | <0.40                                 | <0.40                                 | <0.40                                 |
| Isophorone                   | 0.40                                      | "     | < 0.80                                | < 0.80                                | < 0.80                                |
| 2-Nitrophenol                | 0.14                                      |       | < 0.28                                | <0.28                                 | < 0.28                                |
| 2,4-Dimethylphenol           | 0.17                                      | 71    | < 0.34                                | < 0.34                                | < 0.34                                |
| Bis(2-chloroethoxy)methane   | 0.13                                      | 11    | < 0.26                                | <0.26                                 | < 0.26                                |
| 2,4-Dichlorophenol           | 0.12                                      | 11    | < 0.24                                | <0.24                                 | <0.24                                 |
| 1,2,4-Trichlorobenzene       | 0.20                                      | **    | ≤0:40 R                               | <0.40 P                               | ≈0.40 R                               |
| Naphthalene                  | 0.03                                      | 11    | <0.06                                 | < 0.06                                | < 0.06                                |
| Hexachlorobutadiene          | 0.20                                      | n     | < 0.40                                | <0.40                                 | < 0.40                                |
| 4-Chloro-3-Methylphenol      | 0.14                                      | 0     | < 0.28                                | <0.28                                 | < 0.28                                |
| Hexachlorocyclopentadiene    | 0.20                                      | *1    | < 0.40                                | < 0.40                                | < 0.40                                |
| 2,4,6-Trichlorophenol        | 0.12                                      | 11    | < 0.24                                | < 0.24                                | < 0.24                                |
| 2-Chloronaphthalene          | 0.09                                      | **    | < 0.18                                | < 0.18                                | < 0.18                                |
| Acenaphthylene               | 0.04                                      | 84    | <0.08                                 | < 0.08                                | < 0.08                                |
| Dimethyl phthalate           | 0.11                                      | Ħ     | <0.22                                 | < 0.22                                | < 0.22                                |
| 2,6-Dinitrotoluene           | 0.06                                      | 11    | < 0.12                                | < 0.12                                | < 0.12                                |
| Acenaphthene                 | 0.07                                      | н     | < 0.14                                | < 0.14                                | < 0.14                                |
| 2,4-Dinitrophenol            | 0.48                                      | **    | <0.96                                 | <0.96                                 | <0.96                                 |
| 2,4-Dinitrotoluene           | 0.05                                      | 11    | < 0.10                                | < 0.10                                | < 0.10                                |
| 4-Nitrophenol                | 0.14                                      | **    | <0.28                                 | < 0.28                                | < 0.28                                |
| Fluorene                     | 0.03                                      | **    | < 0.06                                | <0.06                                 | < 0.06                                |
| 4-Chlorophenylphenylether    | 0.09                                      | n     | < 0.18                                | < 0.18                                | < 0.18                                |
| Diethyl phthalate            | 0.11                                      | **    | < 0.22                                | < 0.22                                | < 0.22                                |
| 4,6-Dinitro-2-methylphenol   | 0.15                                      | **    | < 0.30                                | < 0.30                                | < 0.30                                |
| N-Nitrosodiphenylamine       | 0.19                                      | "     | < 0.38                                | < 0.38                                | < 0.38                                |
| 4-Bromophenylphenylether     | 0.03                                      | "     | < 0.06                                | < 0.06                                | < 0.06                                |
| Hexachlorobenzene            | 0.20                                      | "     | < 0.40                                | <0.40                                 | <0.40                                 |
| Pentachlorophenol            | 0.11                                      | 11    | < 0.22                                | <0.22                                 | <0.22                                 |
| Phenanthrene                 | 0.03                                      | *1    | < 0.06                                | < 0.06                                | < 0.06                                |
| Anthracene                   | 0.02                                      | **    | < 0.04                                | < 0.04                                | < 0.04                                |
| Di-n-butyl phthalate         | 0.11                                      | "     | < 0.22                                | < 0.22                                | < 0.22                                |
| Fluoranthene                 | 0.02                                      | **    | < 0.04                                | < 0.04                                | < 0.04                                |
| Pyrene                       | 0.03                                      |       | < 0.06                                | < 0.06                                | < 0.06                                |
| Benzyl butyl phthalate       | 0.06                                      |       | < 0.12                                | < 0.12                                | < 0.12                                |
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|                          | Client ID:<br>Zenon ID:<br>Date Sampled: |       | BH1<br>10-12<br>055067 96<br>96/12/12 | BH2<br>15-17<br>055068 96<br>96/12/12 | BH3<br>13-15<br>055069 96<br>96/12/12 |
|--------------------------|--|-------|---------------------------------------|---------------------------------------|---------------------------------------|
| Component                | MDL                                      | Units | 90/12/12                              | 90/12/12                              | 90/12/12                              |
| 3,3-Dichlorobenzidine    | 0.10                                     | "     | <0.20                                 | <0.20                                 | < 0.20                                |
| Benzo(a)anthracene       | 0.02                                     |       | <0.04                                 | <0.04                                 | <0.04                                 |
| Chrysene                 | 0.03                                     | 11    | < 0.06                                | <0.06                                 | <0.06                                 |
| Bis(2-ethylhexyl)phthala |  | 11    | <0.28                                 | <0.28                                 | <0.28                                 |
| Di-n-octyl phthalate     | 0.11                                     | **    | < 0.22                                | < 0.22                                | < 0.22                                |
| Benzo(b)fluoranthene     | 0.04                                     | *1    | < 0.08                                | < 0.08                                | < 0.08                                |
| Benzo(k)fluoranthene     | 0.04                                     | "     | < 0.08                                | < 0.08                                | <0.08                                 |
| Benzo(a)pyrene           | 0.05                                     | **    | < 0.10                                | < 0.10                                | < 0.10                                |
| Indeno(1,2,3-cd)pyrene   | 0.06                                     | "     | < 0.12                                | < 0.12                                | < 0.12                                |
| Dibenzo(a,h)anthracene   | 0.04                                     | 11    | < 0.08                                | < 0.08                                | < 0.08                                |
| Benzo(ghi)perylene       | 0.04                                     | 11    | < 0.08                                | < 0.08                                | < 0.08                                |
| N-Nitrosodimethylamine   | 1.0                                      | 11    | <2.0                                  | <2.0                                  | <2.0                                  |
| Aniline                  | 0.50                                     | **    | <1.0                                  | <1.0                                  | <1.0                                  |
| Benzyl alcohol           | 0.50                                     | 11    | <1.0                                  | <1.0                                  | <1.0                                  |
| Carbazole                | 0.50                                     | *1    | <1.0                                  | <1.0                                  | <1.0                                  |
| 2-Methylphenol           | 0.50                                     | 11    | <1.0                                  | <1.0                                  | <1.0                                  |
| Benzoic acid             | 0.50                                     |       | <1.0                                  | <1.0                                  | <1.0                                  |
| 4-Chloroaniline          | 0.50                                     | **    | <1.0                                  | <1.0                                  | <1.0                                  |
| 2-Methylnaphthalene      | 0.10                                     | *1    | <0.20                                 | < 0.20                                | < 0.20                                |
| 2,4,5-Trichlorophenol    | 0.10                                     |       | <0.20                                 | < 0.20                                | < 0.20                                |
| 2-Nitroaniline           | 0.50                                     | "     | <1.0                                  | <1.0                                  | <1.0                                  |
| 3-Nitroaniline           | 0.50                                     | **    | <1.0                                  | <1.0                                  | <1.0                                  |
| Dibenzofuran             | 0.50                                     | **    | <1.0                                  | <1.0                                  | <1.0                                  |
| Benzidine                | 0.50                                     | **    | <1.0                                  | <1.0                                  | <1.0                                  |
| 4-Nitroaniline           | 0.50                                     |       | <1.0                                  | <1.0                                  | <1.0                                  |
| Surrogate Recoveries     |  | %     |                                       |                                       |                                       |
| 2-Fluorophenol           |  |       | 39                                    | 43                                    | 43                                    |
| d5-Phenol                |  |       | 66                                    | 65                                    | 68                                    |
| d5-Nitrobenzene          |  |       | 41                                    | 43                                    | 46                                    |
| 2-Fluorobiphenyl         |  |       | 60                                    | 62                                    | 65                                    |
| 2,4,6-Tribromophenol     |  |       | 61                                    | 64                                    | 57                                    |
| d14-p-Terphenyl          |  |       | 72                                    | 70                                    | 72                                    |

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| Batch Code:     | 1220ASA1  |
|-----------------|-----------|
| Mercury         | 055065 96 |
| -               | 055066 96 |
|                 | 055067 96 |
|                 | 055068 96 |
|                 | 055069 96 |
| Date analysed   | 96/12/20  |
| Date prepared   | 96/12/20  |
| Batch Code:     | 1223ASC1  |
| Arsenic         | 055065 96 |
|                 | 055066 96 |
|                 | 055067 96 |
|                 | 055068 96 |
|                 | 055069 96 |
| Date analysed   | 96/12/30  |
| Date prepared   | 96/12/23  |
| Batch Code:     | 1223ASC1  |
| Selenium        | 055065 96 |
|                 | 055066 96 |
|                 | 055067 96 |
|                 | 055068 96 |
|                 | 055069 96 |
| Date analysed   | 96/12/30  |
| Date prepared   | 96/12/23  |
| Batch Code:     | 1223ASC1  |
| Metals          | 055065 96 |
|                 | 055066 96 |
|                 | 055067 96 |
|                 | 055068 96 |
|                 | 055069 96 |
| Date analysed   | 96/12/30  |
| Date prepared   | 96/12/23  |
| Batch Code:     | 1220SM01  |
| Volatiles       | 055065 96 |
|                 | 055066 96 |
|                 | 055067 96 |
|                 | 055068 96 |
|                 | 055069 96 |
| Date analysed   | 96/12/20  |
| Date prepared   | 96/12/20  |
| Batch Code:     | 1217PB02  |
| Semi-Volatiles  | 055065 96 |
| Senn- voiatiles |           |
|                 | 055066 96 |

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Date analysed Date prepared

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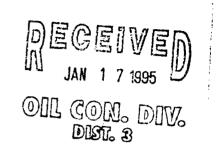
ONE OF THE WILLIAMS COMPANIES ,

L CONSER, SHON SIVISION RECEIPED .

Р.О. Вох разоли Salt Lake City, UT 84158-0900 (2011) 584-7033 25 JAV 18 АЛ 852 P.O. Box 58900

Server a de marce la la marce.

January 13, 1995



Mr. Denny Foust Oil Conservation Division District III 1000 Rio Brazos Road Aztec, New Mexico 87410

Dear Mr. Foust:

At your request, Williams Field Services performed an assessment of the coolant storage tank area at our Milagro Plant. The assessment was performed in order to determine the nature and extent of contamination associated with a release of glycol which occurred on December 3, 1994. The results of the assessment are provided in the attached report. If you have any questions or require additional information, please do not hesitate to contact me at (801) 584-6543.

Sincerely,

Leigh E. Gooding, P.G. Environmental Specialist

ASSESSMENT OF AREA SURROUNDING THE COOLANT STORAGE AT WILLIAMS FIELD SERVICES' MILAGRO PLANT

January 3, 1995

Prepared by:

Williams Field Services Environmental Services Department

#### INTRODUCTION

On Saturday December 3rd, 1994, the Milagro Plant Superintendent notified WFS Environmental Services of the discovery of free glycol in a ditch just outside the south wall of the coolant storage secondary containment located within the facility boundaries of the Milagro Plant. The Milagro Plant is located in San Juan County approximately 3 miles northeast of Bloomfield, New Mexico in Section 12, Township 29N, Range 11W.

#### BACKGROUND

The above mentioned ditch had been cut by an electrical contractor installing conduit between the south wall and a small metal building located only about two feet from the secondary containment. Liquid was later observed to be accumulating in the 18" deep ditch. Further investigation indicated that the liquid appeared to be glycol and water.

Plant personnel subsequently proceeded to take steps to remove the standing glycol from the ditch and to attempt to ascertain the Since the ditch was only one foot south of the secondary source. containment for glycol storage tanks, it was believed that concrete seams (expansion joints) in the concrete wall and floor allowed glycol which had overflowed from the tanks to leak outside the containment designed to prevent same. The overflow of glycol was attributable to two slow leaks discovered in the coolant system which caused the coolant storage tank to overflow. All qlycol which overflowed was thought to be contained within the secondary structure. The "free product" was pumped out of the ditch and from the secondary containment structure by plant personnel within hours of its discovery. It is estimated that approximately 100 gallons total was removed from the ditch and the concrete containment.

After removing all standing glycol from the ditch (apprx. 50 gallons), plant personnel observed no further accumulation of "free product". In order to attempt to delineate the extent of the area impacted by the release of glycol outside the containment area, WFS Environmental Services initiated an assessment action on December 10, 1994.

#### DESCRIPTION OF ASSESSMENT ACTIVITIES

As depicted in the attached site diagram, soil samples were collected from the center point approximately 12" outside of each of the four sides of the containment walls. Samples were collected using a hand driven soil probe from depths between two and three feet. Results obtained following sample analysis are presented in the attached data summary table.

In addition to these samples, two other samples were collected to attempt to determine the areal extent of glycol affected soils. One sample was collected approximately 12' from the southwest corner of the secondary containment in the direction thought to be down gradient at a depth of 3'.

The second sample was collected from a stormwater ditch across a service road running east of the coolant storage area. This sample was collected at a depth of 12". There was no surface staining indicating that any glycol had migrated across the road nor was there evidence of any staining of gravels or road base adjacent to the secondary containment.

#### DISCUSSION OF FINDINGS

After evaluating the analytical results obtained, WFS notes that the only areas impacted by the leaking secondary containment are directly adjacent to the south and east walls. The area outside the south wall was the area where the free product was recovered and that area where glycol impacted soil was obvious. The soils outside the east wall were found to contain glycol at a concentration of 28,000 parts per million.

All other samples showed only trace or non-detectable levels of glycol.

#### CONCLUSION

Based on the results obtained, it appears that the glycol impacted soil is limited to that area immediately south and east of the coolant storage area at the Milagro Plant. WFS plant personnel have already corrected what appears to have been a design flaw in the secondary containment which led to the leakage of free product. This entailed the re-sealing of all concrete seams with material specifically made for this purpose. This action should eliminate future problems of this type and further enhance the integrity of the existing secondary containment.

Because all recoverable "free product" outside the containment has been removed, and because of the limited area which was impacted, WFS believes no further action is necessary. The glycol present in the soils surrounding the coolant storage area will naturally degrade over time. Since the annual precipitation in the area is just over 8 inches, the potential to impact stormwater runoff to any measurable extent is minimal. Likewise, since the depth to ground water is over 100', migration resulting from surface water infiltration to that depth is unlikely.

Plant personnel will continue to check the effectiveness of the secondary containment at the subject site as well as other containments at the facility in accordance with provisions of the Stormwater Pollution Prevention Plan prepared specifically for the Milagro Plant. If at any time, conditions at or around the coolant storage area are perceived to be contributing to a release of any product off-site, then Milagro Plant personnel will initiate immediate corrective action.

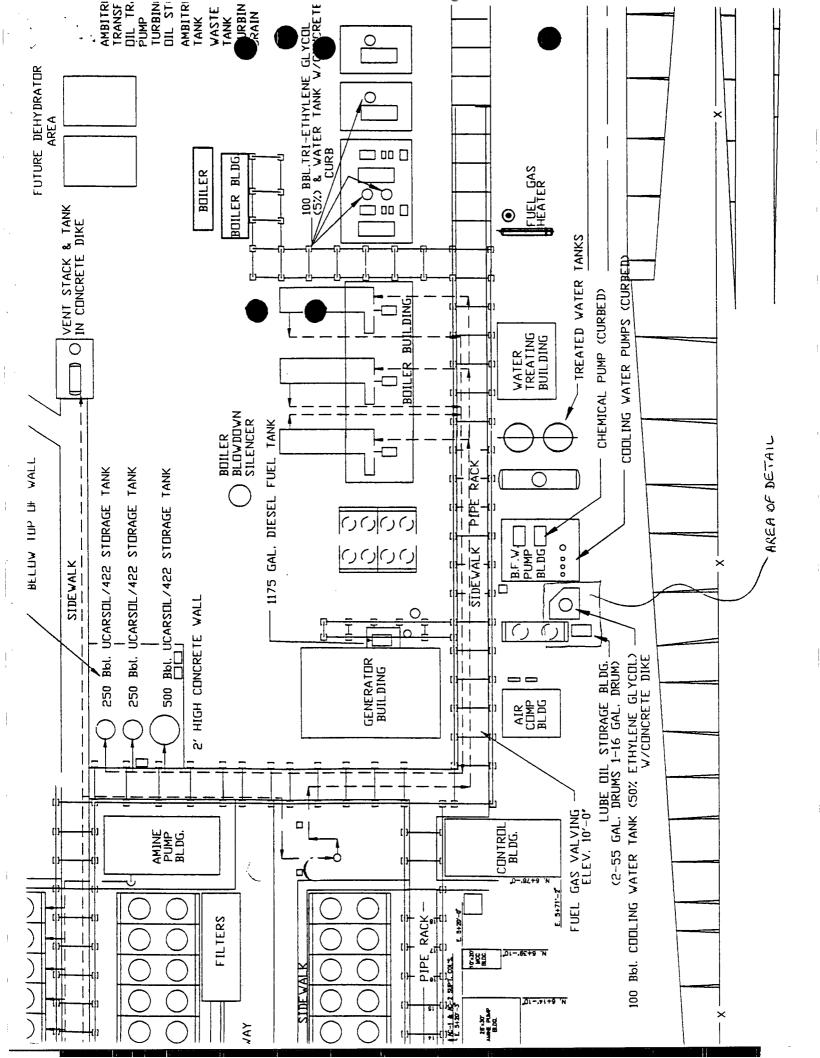
### MILAGRO ASSESSMENT

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#### SUMMARY OF LABORATORY RESULTS

| SAMPLE NUMBER | DETECTION<br>LIMIT (PPM) | CONCENTRATION<br>(PPM) | SAMPLE LOCATION AND DEPTH       |
|---------------|--------------------------|------------------------|---------------------------------|
| MIL-E01       | 10                       | 29,200                 | East Wall 24" Below Surface     |
| MIL-NO1       | 10                       | 27                     | North Wall 30" Below Surface    |
| MIL-SO1       | 10                       | 10,500                 | South Wall 46" Below Surface ** |
| MIL-W01       | 10                       | 91                     | West Wall 36" Below Surface     |
| MIL-SW01      | 10                       | 24                     | Point 12'from SW Corner 36" BS  |
| MIL-DD01      | 10                       | 22                     | Soil from Storm Ditch 12" BS    |

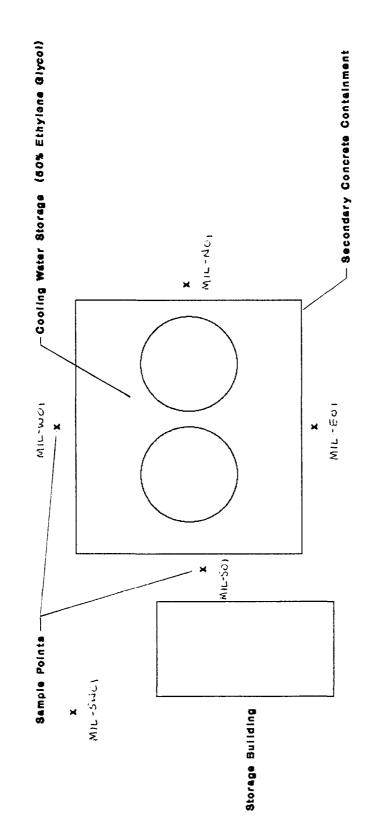
\*\* Note: This sample was collected from within the ditch south of the containment structure. The depth below the bottom of the ditch is approximately 28".



MILAGRO PLANT COOLANT STORAGE AREA

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Service Road

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MIL-DDOI

Stormwater Ditch

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WEST ANALYTICAL LABORATORIES

463 West 3600 South Salt Lake City, Utah 84115

(801) 263-8686 Fax (801) 263-8687

Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

| <u>Analysis Requested:</u><br>Ethylene & Propylene Glycol | Method Ref. Number:<br>Extraction-Direct Injection GC/MS | Date Extracted:<br>December 15, 1994       |
|---|--|--|
| Lab Sample ID, Number:<br>20874-Method Blank              | Field Sample ID, Number:<br>Method Blank                 | <u>Date Analyzed:</u><br>December 16, 1994 |
| Analytical Results<br>Units = mg/kg (ppm)                 | ······································                   |  |
| Compound:   |  | mount<br>etected:                          |
| Ethylene Glycol   | 10.  | <10.                                       |
| Propylene Glycol  | 10.  | <10.                                       |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

| Released by: | - Or           | ( some   | 土 |       |
|--------------|----------------|----------|---|-------|
| ·            | Laboratory Sup | pervisor | ) | <br>- |

Report Date 12/20/94

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Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

|  | nod Ref. Number:<br>action-Direct Injection GC/MS | Date Extracted:<br>December 15, 1994 |
|--|---|--------------------------------------|
|--|---|--------------------------------------|

Lab Sample ID. Number: 20874-01

Field Sample ID, Number:Date Analyzed:Milagro Plant-SJADecember 16, 1994MIL-EO1/Soil From E. Side of CoolantStorage 24" B.S.

### Analytical Results Units = mg/kg (ppm)

| (801) 263-8686<br>Fax (801) 263-8687 | Compound:        | Detection<br>Limit: | Amount<br>Detected: |
|--------------------------------------|------------------|---------------------|---------------------|
|                                      | Ethylene Glycol  | 10.                 | 28,000.             |
|                                      | Propylene Glycol | 10.                 | 1,200.              |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

| Released by: | - Ore         | CA        | ,t |
|--------------|---------------|-----------|----|
| ·            | Laboratory Su | pervisor/ | )  |

Report Date 12/20/94

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463 West 3600 South Salt Lake City, Utah 84115



Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples

Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

| Analysis Requested:Method Ref. NoEthylene & Propylene GlycolExtraction-Dire | Imber:Date Extracted:ct Injection GC/MSDecember 15, 1994 |
|---|--|
|---|--|

Lab Sample ID. Number: 20874-02

Field Sample ID. Number:Date Analyzed:Milagro Plant-SJADecember 16, 1994MIL-NO1/Soil From N. Side of CoolantStorage 30" B.S.

## Analytical Results Units = mg/kg (ppm)

| Compound:        | Detection<br>Limit: | Amount<br>Detected: |
|------------------|---------------------|---------------------|
| Ethylene Glycol  | 10.                 | 15.                 |
| Propylene Glycol | 10.                 | 12.                 |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

| Released by: | Doe count             |  |
|--------------|-----------------------|--|
| -            | Laboratory Supervisor |  |

Report Date 12/20/94

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463 West 3600 South Salt Lake City, Utah 84115

(801) 263-8686 Fax (801) 263-8687



463 West 3600 South Salt Lake City, Utah

(801) 263-8686 Fax (801) 263-8687

84115

**ORGANIC ANALYSIS REPORT** 

Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

| Analysis Requested:<br>Ethylene & Propylene Glycol   | Method Ref. Number:<br>Extraction-Direct Injection GC/MS | <u>Date Extracted:</u><br>December 15, 1994 |
|--|--|---|
| <u>Lab Sample ID. Number</u> :<br>20874-Method Blank | Field Sample ID, Number:<br>Method Blank                 | <u>Date Analyzed:</u><br>December 15, 1994  |
| Analytical Results<br>Units = mg/kg (ppm)            |  |   |
| Compound:  |  | mount<br>tected:                            |
| Ethylene Glycol                                      | 10.  | <10.  |
| Propylene Glycol                                     | 10.  | <10.  |
|  |  |   |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

Released by: Laboratory Supervisor

Report Date 12/20/94

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# **ORGANIC ANALYSIS REPORT**

Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

| Analysis Requested:         | Method Ref. Number:               | <u>Date Extracted:</u> |
|-----------------------------|-----------------------------------|------------------------|
| Ethylene & Propylene Glycol | Extraction-Direct Injection GC/MS | December 15, 1994      |

Lab Sample ID. Number: 20874-03

Field Sample ID. Number:Date Analyzed:Milagro Plant-SJADecember 15, 1994MIL-SO1/Soil From S. Side of CoolantStorage 46" B.S.

## Analytical Results Units = mg/kg (ppm)

| Compound:        | Detection<br>Limit: | Amount<br>Detected: |
|------------------|---------------------|---------------------|
| Ethylene Glycol  | 10.                 | 9,400.              |
| Propylene Glycol | 10.                 | 1,100.              |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

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Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

Analysis Requested:Method Ref. Number:Date Extracted:Ethylene & Propylene GlycolExtraction-Direct Injection GC/MSDecember 15, 1994

Lab Sample ID, Number: 20874-04

Field Sample ID, Number:Date Analyzed:Milagro Plant-SJADecember 15, 1994MIL-WO1/Soil From W. Side of CoolantStorage 36" B.S.

### Analytical Results Units = mg/kg (ppm)

| Compound:        | Detection<br>Limit: | Amount<br>Detected: |
|------------------|---------------------|---------------------|
| Ethylene Glycol  | 10.                 | 79.                 |
| Propylene Glycol | 10.                 | 12.                 |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

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# **ORGANIC ANALYSIS REPORT**

Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples

Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

Method Ref. Number: Extraction-Direct Injection GC/MS

iction-Direct Injection GC/MS

Date Analyzed: December 15, 1994

December 15, 1994

Date Extracted:

<u>Lab Sample ID, Number</u>: 20874-05

Ethylene & Propylene Glycol

Analysis Requested:

Field Sample ID. Number:DatMilagro Plant-SJADecMIL-SWO1/Soil From Point 12' fromSW Corner of Containment @ 36" B.S.

### Analytical Results Units = mg/kg (ppm)

| Compound:        | Detection<br>Limit: | Amount<br>Detected: |
|------------------|---------------------|---------------------|
| Ethylene Glycol  | 10.                 | 12.                 |
| Propylene Glycol | 10.                 | 12.                 |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

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84115

# **ORGANIC ANALYSIS REPORT**

Client: Williams Field Services Date Sampled: December 10, 1994 Date Received: December 12, 1994 Set Description: Six Waste Samples Set Identification #: 20874 Contact: Mark Harvey Received By: Sherlyn Lewis

Analysis Requested:Method Ref. Number:Date E:Ethylene & Propylene GlycolExtraction-Direct Injection GC/MSDecem

Lab Sample ID. Number: 20874-06

<u>Field Sample ID. Number</u>: Milagro Plant-SJA MIL-DDO1/Soil From Drainage E. of Coolant Containment @ 16" Date Extracted: December 15, 1994

Date Analyzed: December 15, 1994

## Analytical Results Units = mg/kg (ppm)

| Compound:        | Detection<br>Limit: | Amount<br>Detected: |
|------------------|---------------------|---------------------|
| Ethylene Glycol  | 10.                 | 11.                 |
| Propylene Glycol | 10.                 | 11.                 |

< Value = None detected above the specified method detection limit, or a value that reflects a reasonable limit due to interferences.

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