

SECOND INTERIM REMEDIAL MEASURES WORK PLAN FORMER SIGNORE FACILITY ELLICOTTVILLE, NEW YORK BROWNFIELD CLEANUP PROGRAM SITE NO. C905034

# **Prepared For:**

New York State Department of Environmental Conservation Region 9 Buffalo, New York

# **Prepared By:**

GZA GeoEnvironmental of New York Buffalo, New York



July 2013

File No: 21.0056367.50

Copyright© 2013 GZA GeoEnvironmental of New York

# SECOND INTERIM REMEDIAL MEASURES WORK PLAN FORMER SIGNORE FACILITY ELLICOTTVILLE, NEW YORK BROWNFIELD CLEANUP PROGRAM SITE NO. C905034

TABLE OF CONTENTS

|                         | Page  | <u> </u> |
|-------------------------|---|----------|
| 1.0 INTROD              | OUCTION1  | Ĺ        |
| 1.2 PROJE<br>1.3 PROJE  | OSE AND OBJECTIVE   | 1<br>1   |
|                         | ECT MANAGEMENT AND ORGANIZATION4  |          |
|                         | ersonnel4 pecific Tasks and Services4   |          |
| 2.0 DESCRI              | PTION OF IRM ACTIVITIES5  | 5        |
| 2.1.1 Site              | RAL FIELD ACTIVITIES  | 5        |
|                         | obilization   |          |
|                         | URVEILLANCE AND MONITORING  |          |
| 2.3 SOIL I<br>2.4 ELECT | EXCAVATIONS6<br>TRON DONOR COMPOUND INJECTIONS & GROUNDWATER SAMPLING 8<br>RONMENTAL ANALYTICAL TESTING PROGRAM11 | 5<br>3   |
| 3.0 DATA D              | OCUMENTATION11  | ĺ        |
| 4.0 IRM AC              | CTIVITIES REPORTING12   | 2        |
| 5.0 QUALIT              | Y ASSURANCE/QUALITY CONTROL12   | 2        |
| 6.0 HEALTH              | H AND SAFETY PROTOCOLS  | 3        |
| 7.0 CITIZEN             | PARTICIPATION13   | 3        |
| 8.0 SCHEDU              | JLE   | 3        |
| TABLES                  |   |          |
| Table 1                 | Estimated IRM Analytical Testing Program Summary  |          |
| FIGURES                 |   |          |
| Figure 1                | Locus Plan  |          |
| Figure 2                | Site Plan   |          |
| Figure 3                | Chlorinate VOC Concentration Isopleth   |          |
| Figure 4                | AOC-2, AOC-3 and Injection Location Plan  |          |

# SECOND INTERIM REMEDIAL MEASURES WORK PLAN FORMER SIGNORE FACILITY ELLICOTTVILLE, NEW YORK BROWNFIELD CLEANUP PROGRAM SITE NO. C905034 TABLE OF CONTENTS (CONT'D)

# **APPENDICES**

| Appendix A | Table 3 Phase II Soil Results Summary Table |
|------------|---|
| Appendix B | SVOC and METALS Results From AOC-2          |
| Appendix C | Electron Donor Compound Material MSDS       |

# 1.0 INTRODUCTION

# 1.1 PURPOSE AND OBJECTIVE

This Second<sup>1</sup> Interim Remedial Measures (IRM) Work Plan has been developed by GZA GeoEnvironmental of New York (GZA), on behalf of Iskalo Ellicottville Holdings, LLC (Iskalo), for IRM activities associated with the Former Signore Facility Brownfield Cleanup Program (BCP) Site No. C905034 located at 55-57 Jefferson Street, Ellicottville, New York (Signore BCP Site, see Figures 1 and 2). Specifically, this work plan has been developed to address the remaining soil contamination associated with Area of Concern-2 (AOC-2) and AOC-3; and complete a pilot test to assess chlorinated volatile organic compound (cVOC) groundwater contamination treatment.

The work described in this IRM Work Plan is being done under a New York State Department of Environmental Conservation (NYSDEC) BCP Agreement. This IRM Work Plan presents the project scope, objectives, planned activities, sampling procedures and reporting requirements.

# 1.2 PROJECT BACKGROUND

The Former Signore Facility is located at 55-57 Jefferson Street in the Village of Ellicottville, Cattaraugus County, New York. Ellicottville is located approximately 60 miles south of Buffalo, New York, and is a popular ski-resort area. General adjoining land uses are residential and recreational. The entire property is approximately 55 acres of which 8.43 acres (Signore BCP Area, see Figure 2) are occupied by the concrete slab of the former Signore building (168,000 square feet), other ancillary buildings and parking areas. The remaining acreage is vacant, undeveloped land. The property consists of approximately 21 acres of "flat land" area and about 34 acres of hill side.

The property has been used for manufacturing purposes for over 50 years. It is reported that a tool and die operation occupied a garage associated with the residential dwelling that was formerly present at the property. The Signore BCP Area was primarily used for the manufacturing of metal products. The existing Signore building has undergone various expansions since 1952. The actual development date for the property is unknown, but occurred sometime between the 1940s and 1952 as the property was identified as vacant woodland between 1922 and 1939.

The property is listed on the NYSDEC State Superfund Program as Site number 905023. In 1986, the Signore facility undertook a soil and groundwater sampling program which identified low concentrations of volatile organic compounds (VOCs) at the Site. Both downgradient public and private drinking water wells were affected. The contamination was attributed to spills,

\_

<sup>&</sup>lt;sup>1</sup> The original IRM Work Plan entitled "Interim Remedial Measures Work Plan, Former Signore Facility, Ellicottville, New York, Brownfield Cleanup Program, Site No. C905034" dated July 2011, prepared by GZA, addressed the areas of AOC-1 and the majority of AOC-2.

leakage and other plant operations.

In August 1989, Signore entered into an Administrative Order on Consent #89-258-89-03 to perform a Remedial Investigation/Feasibility Study (RI/FS) at the Site and three Interim Remedial Measures (IRMs). The three IRMs included the following.

- 1. Installation of an interceptor well upgradient of the Town drinking water well;
- 2. Connection of 34 residential properties to the municipal water supply source;
- 3. Installation of an interceptor well on a downgradient portion of the Signore property.

The IRM activities were completed and in operation by January 1992. The contaminant of concern was identified as trichloroethene (TCE) and trichloroethane (TCA). Additional volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), metals, or polychlorinated biphenyls (PCBs) were not identified on the Site during the previous RI.

In 1993, the Site was reclassified from a Class #2 to a Class #4 Site, as it has been properly closed. However, NYSDEC requires groundwater monitoring at on and off-site locations on a semi-annual basis, which Iskalo is continuing to perform. Long term monitoring data has shown a general decrease in site contaminants and off-site migration.

In 2002, the on-Site interceptor well and the Town Well interceptor well were shut down, as approved by NYSDEC, due to long-term sampling results at both wells indicating either non-detect concentrations or levels below State drinking water standards.

In October 2007, GZA completed a Phase II ESA at the Site as part of due diligence services for Iskalo. GZA's work included observing soil probes at 29 locations and test pit excavations at eight locations. During the Phase II activities, VOC contamination and separate phase petroleum (SPP) product were identified impacting soil and groundwater at the Site. Three areas of concern (AOC) were identified where the contaminant concentrations were greater than the NYSDEC Unrestricted Soil Cleanup Objectives (6 NYCRR Part 375<sup>2</sup> criteria) [Table 3 from the Phase II ESA, which summarized the results of the soil samples has been included in Appendix A for reference.]

A brief description of each AOC follows below. AOC-1 and the majority of AOC-2 were addressed under the previous original NYSDEC-approved IRM work plan. The activities associated with this work were completed between October 2011 and December 2011. AOC-3 was located beneath the existing building during the time of the first IRM which has since been demolished.

<sup>&</sup>lt;sup>2</sup> 6 New York Code Rules and Regulation (6 NYCRR) Part 375 Environmental Remediation Programs, effective December 14, 2006 (Part 375).

- 1. AOC-1 Petroleum underground storage tanks (UST) Area Three 1,000-gallon USTs, located on the eastern portion of the Site, were closed in-place in December 1986. SPP product and petroleum impacted soil was identified during test pit completion. GZA contacted NYSDEC and Spill #707350 was assigned to the Site on October 3, 2007. This area was addressed as part of the IRM activities completed in 2011.
- 2. AOC-2 One 1,000-gallon UST Area The historic contents of a UST identified on the southwest side of the main building are unknown. The UST was reportedly closed in the late 1980s. The majority of this area was addressed as part of the IRM activities completed in 2011. Some contamination remains on the north wall, which was located along the building foundation and under the building at the time the work was completed. The building has since been demolished.
- 3. AOC-3 Paint Kitchen Area VOC impacted soil was identified under the main building identified as the paint kitchen and spray booth area. Additionally, a former septic system was also present in the area. "Product" was identified during the soil probe investigation. This area is located under the remaining concrete slab and is to be addressed as part of the IRM activities discussed in this IRM Work Plan.

A Supplemental Remedial Investigation (SRI) was completed between January 2012 and January 2013. The activities included the following.

- Off-site soil vapor intrusion assessment of nine homes;
- Completion of ten (10) on-site test pits;
- Completion of 21 on-site soil probes;
- Collection and analysis of four (4) on-site surface soil samples;
- Collection and analysis of 21 soil samples from the 21 soil probes; and
- Collection of 19 groundwater samples from the 14 new microwells installed as part of the SRI and five (5) existing wells.

Groundwater contamination was further defined within the Signore BCP Site boundaries, during the SRI activities completed in Fall 2012. The results indicated cVOCs are present on-site and at concentrations greater than the Class GA criteria. Figure 3 is a total cVOC isopleth map generated using data generated as part of the SRI and previous Phase II ESA.

The petroleum-based VOC impacts appear to be associated with residual petroleum source material in the soil and do not appear to be migrating from the source areas (AOCs -1, -2, and -3). The cVOCs in the groundwater do not appear to be related to former AOC-1 or AOC-3, but are likely related to other historical on-Site activities.

Therefore, the purpose of the pilot test, as discussed later in this work plan, is to assess a treatment technology to treat the cVOC contamination identified in the Signore BCP Site groundwater.

# 1.3 PROJECT DESCRIPTION

The IRM activities will involve the removal of impacted soils located in the vicinity of AOC-2 and AOC-3. See Figure 4 for the approximate locations of impacted soil to be addressed at AOC-2 and AOC-3 (shown in blue). The green boxed areas shown on Figure 4 represent the approximate areas where the electron donor compound (EDC) material will be injected in the cVOC impacted groundwater as part of the pilot test.

# 1.4 PROJECT MANAGEMENT AND ORGANIZATION

# 1.4.1 Personnel

The general responsibilities of key project personnel are listed below.

*NYSDEC Project Manager* - Chad Staniszewski will have the responsibility for regulatory oversight for the work associated with BCP Site No. C905034.

*Iskalo Development Project Manager* – Paul B. Iskalo will have the responsibility for implementing the project and has the authority to commit funding necessary to meet the objectives and requirements.

GZA Project Manager - Christopher Boron will be responsible for managing the implementation of the activities associated with the BCP investigation, remediation and coordinating the collection of data during the project. The Project Manager is responsible for technical quality control and project oversight.

GZA Quality Assurance (QA) Officer – Daniel Troy P.E., will report to the Project Manager and will be responsible for ensuring that QA/QC procedures are being followed. The QA Officer will be responsible for overseeing the review of field and laboratory data.

The QA Officer will monitor the performance of the laboratory to verify that the Data Quality Objectives for the project are met.

GZA Field QA Officer – Thomas Bohlen will be responsible for the overall operation of the field team and reports directly to the Project Manager.

#### 1.4.2 Specific Tasks and Services

Subcontractor specialists will be utilized for services relating to the IRM and are as follows.

<u>Matrix Environmental Technologies, Inc.</u> – Soil excavation, contaminated soil removal, soil disposal.

<u>Spectrum Analytical Services</u> - Laboratory/analytical services

<u>Data Validation Services</u> – Data usability summary report of laboratory data.

# 2.0 DESCRIPTION OF IRM FIELD ACTIVITIES

# 2.1 GENERAL FIELD ACTIVITIES

General field activities include site meetings, mobilization, implementing the health and safety plan, soil excavations, analytical testing, equipment decontamination, handling of investigation wastes, and mixing and injection of EDC material.

# 2.1.1 Site Meeting

A Site "kick-off" meeting will be held with Iskalo, GZA and the earth work subcontractor prior to initiating field work activities. The purpose of the meeting will be to orient field team members, Iskalo staff and subcontractors with the Site, project personnel, Site background, scope of work, potential dangers, health and safety requirements, site-specific security and safety protocols, emergency contingencies and other field procedures. NYSDEC staff are welcome to attend and will be notified at least seven (7) days in advance of the meeting.

#### 2.1.2 Mobilization

Following approval of the IRM Work Plan by NYSDEC, the Underground Facilities Protection Organization (UFPO) will be contacted at 1-800-962-7962 to clear exploration locations. Utility clearance will require three working days by UFPO. GZA and its subcontractors will then mobilize necessary materials and equipment to the Site.

# 2.1.3 Health and Safety

It is anticipated that the work to be completed at the Site will be done at level D personal protection. Should health and safety monitoring during field activities warrant an upgrade to level C protection, work will stop and Site conditions will be re-evaluated prior to further investigation activities. See Section 6.0 for additional information on Health and Safety.

# 2.2 COMMUNITY AIR MONITORING PLAN

During the IRM excavation activities, a Community Air Monitoring Plan will be implemented which will include screening for total volatile organics and particulate levels at the perimeter of the work area for health and safety concerns. Section 6.0 of this document references the Health and Safety Plan which will be used for this work [Section 5.2 and Appendix C of the Health and Safety Plan discuss the New York State Department of Health Generic Community Air Monitoring Plan that will be implemented.]

Organic vapors will be monitored with a portable organic vapor meter (OVM) equipped with a photoionization detector (PID) using a 11.7 electron volt (eV) bulb. The particulates will be monitored using equipment that is capable of measuring particle sizes greater than 10-micrometers (PM-10) and can integrate measurements over a 15-minute time frame. The equipment will also have an audible alarm indicating an exceedance of the action level.

Additional details on the total organic and particulate monitoring are presented in the Site-specific Health and Safety Plan (see Section 6.0).

# 2.3 SOIL EXCAVATIONS

# AOC 2 - Impacted Soil Area

A concrete UST and associated impacted soil were removed as part of the IRM activities in 2011. However, due to the presence of the building at the time of the work, some impacted material was left along the building foundation and likely extend under the floor slab. The building has been demolished. Figure 4 identifies the approximate location the remaining impacted soil. It is estimated that approximately 150 cubic yards or about 250 tons of impacted soil remain.

# AOC 3 - Impacted Soil Area

Petroleum-based VOC impacted soil was identified in the area under the main building identified as the paint kitchen and spray booth area. Additionally, a former septic system was also present in the area. Groundwater sample results from downgradient or sampling locations adjacent to AOC-3 did not indicate the presence of petroleum-VOCs. It appears the petroleum-based VOC groundwater contamination is isolated to this area. The estimated soil contamination area is approximately 60 feet by 115 feet (6,900 square feet), with impacts extending to the water table at approximately 12 feet bgs. Figure 4 identifies the approximate location of the remaining impacted soil. It is estimated that approximately 3,000 cubic yards or about 4,800 tons of impacted soil remain.

# Septic Tank Removals

Two (2) closed-in-place septic tanks are located in the central portion of the Signore BCP Site under the concrete slab (see Figure 4). These tanks were reported cleaned and filled with concrete. The tanks will be broken up for removal and disposed of as solid waste. If impacted soil is encountered in the area of the septic tanks during removal, the impacted soil will also be removed from the subsurface for off-site disposal. Soil samples will be collected from the sidewalls of the excavation for VOC analysis.

# General UST and Soil Removal Procedures

The following are the general impacted soil removal procedures that will be followed for the impacted soil areas.

- Waste characterization samples will be collected from AOC-2 and AOC-3. Based on the estimated volume of soil to be removed, five (5) confirmation soil samples (1 from AOC-2 and 4 from AOC-3) will be sent to the laboratory for analysis.
- Prior to the start of the subsurface soil work, a decontamination pad will be built to allow equipment used during the excavation activities to be decontaminated. The pad will be constructed on a stable onsite surface using a minimum of 6-mil plastic sheeting and allow water generated during the decontamination processes to be contained and transferred to 55-gallon drums for characterization and proper disposal. Upon completion of the excavation work, the decontamination pad will be disposed of with impacted soils stockpiled for off-site disposal at a permitted disposal facility.
- Overburden soil excavated to facilitate the removal of the impacted material will be field screened to determine if the soil is potentially impacted. Non-impacted soil (those that do not appear to be visually impacted and registering 10 ppm or less on an OVM during field screening) will be excavated and stockpiled, on 6-mil thick plastic sheeting and covered with 6-mil plastic sheeting, for eventual reuse onsite. Representative soil samples from the soil stockpile will be sampled for VOC analysis prior to reuse.
- Tank contents (i.e., liquids and tank sludge, if any) will be removed, containerized and characterized for proper off-site disposal. The tank contents will be properly disposed of by the contractor and disposal documentation will be provided.
- The Contractor will breakup and remove the septic tanks from the excavation. The concrete will be properly disposed of by the contractor and disposal documentation will be provided.
- Soils determined to be impacted, within AOC-2, AOC-3 and septic tank areas, based on visual observations and field screening (registering 10 ppm or greater), will be excavated and loaded for off-site disposal. The soil will be loaded into dump trucks or dump trailers, covered, and transported by a licensed hauler to a permitted solid waste landfill for proper disposal. Disposal documentation will be provided.
- Post-excavation soil samples will be collected from the side walls and bottom of each excavation in accordance with NYSDEC guidelines to confirm that the remaining soil meets, at a minimum, the Part 375 Commercial SCOs. One sidewall sample will be collected from each excavation sidewall that is less than 30 linear feet. If an excavation sidewall exceeds 30 linear feet, then one sample will be collected for every 30 linear feet. A minimum of one bottom sample will be collected from each UST excavation area. One sample will be collected for every 900 square feet of bottom excavation area.

- Confirmatory samples will be analyzed for Target Compound List (TCL) VOCs via EPA Method 8260. Detection limits of the sample analysis will be below the Part 375 Commercial SCOs. [VOCs are the compounds of concern associated with AOC-2 and AOC-3. Soil sample results from the previous IRM activities did not identify SVOCs as a concern for AOC-2. Two (2) of the waste characterization composite samples collected from AOC-3 (WC-2 and WC-3 as shown on Figure 4) were also analyzed for SVOCs (Base-Neutrals) via EPA Method 8270D and Target Analyte List (TAL) metals via EPA Methods 6010C/7471B. Waste characterization sample WC-2 was collected from impacted soil encountered from 10 to 12 feet below ground surface and WC-3 was collected from impacted soil encountered from 8 to 16 feet below ground surface. No SVOCs were detected above method detection limits and the metal analytes detected were below the Part 375 Unrestricted SCOs. Appendix B contains the laboratory report.]
- The excavation areas will be barricaded to keep personnel away from the excavation while awaiting analytical results and prior to backfilling. If post-excavation soil samples indicate that impacted soil remains, it is anticipated that additional soil will be excavated for off-Site disposal.
- If groundwater is encountered within an excavation, a sample will be collected for TCL VOC analysis.
- After the post-excavation soil samples are reviewed and evaluated to be acceptable, and prior to backfilling, PVC injection piping may be installed along the bottom of the excavation to facilitate additive injections at a later time, especially if the excavation and contamination extend down to or below the groundwater table.
- Photographic documentation of the IRM activities will be done and included in the IRM Report (see Section 4.0).
- Upon soil excavation completion, equipment will be decontaminated prior to being removed from the Site at the decontamination pad location.
- Suitable backfill material will be placed and compacted in lifts within the excavation areas. Backfill brought to the Signore BCP Site will meet the requirements outlined in Part 375-6.7(d) and DER-103 Section 5.4(e).

# 2.4 ELECTRON DONOR COMPOUND INJECTIONS & GROUNDWATER SAMPLING

The groundwater data indicates that the cVOC plume (greater than 200 ppb) is originating in the vicinity of the former septic tanks in the central portion of the Signore BCP Site. Prior to

\_\_\_

<sup>3</sup> NYSDEC, Division of Environmental Remediation (DER), DER-10 / Technical Guidance for Site Investigation and Remediation, issued May 3, 2010.

implementing a full-scale in-situ groundwater treatment program, a pilot test will be completed, as discussed below, to assess the effectiveness of the remedial alternative and collect pre-design data.

# Electron Donor Compound Injections

The pilot test will consist of the injection of approximately 2,000 pounds of EDC material in the vicinity of SP-3 and the 200 ppb contour, 500 pounds of EDC in the vicinity of 100 ppb contour near SP-32, and monitoring the groundwater conditions in the areas of the injections as shown on Figure 4. The EDC material is based on food-grade vegetable oils and surfactants. A material safety data sheet (MSDS) for the material has been included in Appendix C.

The EDC material will enhance the anaerobic breakdown of the "parent" cVOCs present at the Signore BCP Site (trichloroethene (TCE), tetrachloroethene (PCE) and 1,1,1 trichloroethane (111-TCA)) via reductive dehalogenation to the "daughter" breakdown product (cisdichloroethene (cis-DCE) and vinyl chloride (VC)), which readily degrade under aerobic conditions.

Reductive dehalogenation is defined as the biologically-mediated replacement of chlorine (as chloride) on a chlorinated organic compound such as PCE or TCE with elemental hydrogen in the presence of a suitable electron donor causing a transformation of the contaminant to a less chlorinated product. An electron donor is defined as a compound capable of supplying electrons during oxidation-reduction reactions. Microorganisms obtain energy by transferring electrons from electron donors such as organic compounds or by the reduction of inorganic compounds to a terminal electron acceptor (TEA). Electron donors are chemically-reduced materials such as fuel hydrocarbons or naturally-occurring organic carbon, which become chemically oxidized during transformation.

For example, reductive dehalogenation of chlorinated VOCs typically occurs sequentially from PCE to TCE, TCE to 1,2-DCE, 1,2-DCE to VC, and VC to ethene and chloride, and ultimately ethene to carbon dioxide and water.

Currently, about 2,500 pounds of EDC<sup>4</sup> material is stored at the Signore BCP Site. The material will be mixed into slurry and injected into the subsurface groundwater. The following injection methodology is proposed for the two areas identified by nearby soil probe location.

#### SP-3 Area

A direct push soil probe unit will be used to advance the probe of the injection equipment. Injection points will be completed in an approximate 15-foot horizontal spacing, over a 40 foot by 40 foot area, for a total of ten (10) injection points.

<sup>4</sup> EDC is manufactures by EcoCycle of Toyama, Japn to be used as a remedial additive to drive the reductive dechlorination of cVOCs.

Injections will occur in two (2) intervals below the groundwater table at each injection location. The depth of the injection intervals will vary from 12 to 20 feet below ground surface (bgs). Injections will be completed by advancing the probe to the first depth injection interval, injecting the required quantity of EDC material, and then advancing the probe deeper to the next injection interval and injecting the required quantity of EDC material using a top-down approach.

A total of 1,950 pounds of material will be used in this injection area. The 1,950 pounds of material will be mixed with about 2,000 gallons of water to create an injectable slurry. The EDC material and water will be mixed on-site in tanks until the EDC material has dissolved into solution. Once dissolved, the slurry will be injected into the subsurface under pressure using a grout pump or equivalent. The injection rate will be limited to the rate at which the formation will accept the slurry material. The injection pressure will be regulated and monitored to avoid "blow-back" up the sides of the injection rods and up previous completed injection points.

The injection quantities are approximately 195 pounds of EDC material per injection location (97.5 pounds per injection interval) using approximately 200 gallons of water per injection location (100 gallons per injection interval).

# SP-32 Area

A direct push soil probe unit will be used to advance the probe of the injection equipment. Injection points will be completed in an approximate 10-foot horizontal spacing, over a 20 foot by 20 foot area, for a total of five (5) injection points.

Injections will occur in two (2) depth intervals below the groundwater table, at each injection location, ranging from approximately 12 to 20 feet below ground surface (bgs). Injection will be completed by advancing the probe to the first depth injection interval, injecting the required quantity of EDC material, and then advancing the probe deeper to the next injection interval and injecting the required quantity of EDC material using a top-down approach.

A total of 500 pounds of material will be used in this injection area. The 500 pounds of material will be mixed with about 550 gallons of water to create an injectable slurry. The EDC material and water will be mixed on-site in tanks until the EDC material has dissolved into solution. Once dissolved, the slurry will be injected into the subsurface under pressure. The injection rate will be limited to the rate at which the formation will accept the slurry material. The injection pressure will be regulated and monitored to avoid "blow-back" up the sides of the injection rods and up previous completed injection points.

The injection quantities are approximately 100 pounds of EDC material per injection location (50 pounds per injection interval) using approximately 110 gallons of water per injection location (55 gallons per injection interval).

# **Groundwater Sampling**

Groundwater samples will be collected from select monitoring locations within the Signore BCP Site to assess the effectiveness of the EDC injections. Two (2) sampling events will be conducted:

- 1) within 3 months (Fall 2013) of the pilot test injections; and
- 2) 9 to 12 months (late Spring 2014) after the pilot test injections.

Groundwater samples will be collected from approximately six (6) locations (EW-1.25, SP-32, SP-37, SP-38, SP-43, and SP-45) shown in red on Figure 4. The following methodologies will be used for the groundwater sample analysis.

VOC: SW-846, 8260B

Methane: RSK175
TOC: EPA 9060
Chloride: EPA 300
Nitrate: EPA 335.2
Sulfate: EPA 300
Ethane: RSK-175
Ethene: RSK-175

Dissolved Iron: SW-846, 6010B Dissolved Manganese: SW-846, 6010B

A water quality meter and flow-through cell will be used to collect field measurements for pH, specific conductance, dissolved oxygen (DO), turbidity, oxygen reduction potential (ORP) and temperature. Disposable polyethylene tubing and a variable speed low-flow sampling pump will be utilized during the sampling events.

# 2.5 ENVIRONMENTAL ANALYTICAL TESTING PROGRAM

The estimated soil and groundwater environmental testing program is summarized in Table 1. The actual number of soil samples will vary based on the size of excavations, observations and engineering judgment. The samples collected as part of this IRM will be subject to analytical testing methodologies that follow NYSDEC Analytical Service Protocol (ASP) Category B deliverables and allow for the development of a data usability summary report (DUSR). Further information regarding sampling and testing methodologies can be found in the Quality Assurance Project Plan QAPP (see Section 5.0).

#### 3.0 DATA DOCUMENTATION

Field notes will be recorded during the IRM work and become part of the project file. The daily field summaries will include the following daily information for the IRM activities:

- Date:
- Meteorological conditions (temperature, wind, precipitation);
- Site conditions (e.g., dry, damp, dusty, etc.);
- Identification of crew members (GZA and subcontractor present) and other personnel (e.g., agency or site owner) present;
- Description of field activities;
- Location(s) where work is performed;
- Sampled collected;
- Problems encountered and corrective actions taken;
- Records of field measurements or descriptions recorded;
- Injection locations and volumes; and
- Notice of modifications to the scope of work.

Photographic documentation of the IRM activities will be done.

#### 4.0 IRM ACTIVITIES REPORTING

The IRM Activities will be included in the Supplemental Remedial Investigation and Alternative Analysis Report. It will summarize the work conducted as part of the impacted soil removals at AOC-2, AOC-3 and the EDC injections. The information will include the following.

- summary of the activities completed as part of the IRM Work;
- present the analytical data from the confirmatory samples collected. Tables containing the analytical results will identify laboratory qualifiers assigned to the data and will identify the detection limits for non-detected compounds (e.g., < 0.5);
- provide figures showing the size and location of IRM activities along with confirmatory sample locations and EDC injection locations;
- provide pertinent photographic documentation of the activities completed; and
- present the disposal documentation of the various material generated for disposal.

# 5.0 QUALITY ASSURANCE/QUALITY CONTROL

The Quality Assurance Project Plan (QAPP) to be used for the Former Signore Facility IRM activities is the "Quality Assurance Project Plan, Former Signore Facility, Ellicottville, New York, Brownfield Cleanup Program, Site No. C905034" dated May 2011. The QAPP presents the sampling procedures, analytical methods and QA/QC procedures associated with the activities planned for the BCP Site. Protocols for sample collection, sample handling and storage, Chain of Custody procedures, and laboratory and field analyses are described or specifically referenced to related investigation documents.

# 6.0 HEALTH AND SAFETY PROTOCOLS

The health and safety protocols to be used for the Former Signore Facility IRM activities are in the "Health and Safety Plan, Former Signore Facility, Ellicottville, New York, Brownfield Cleanup Program, Site No. C905034" dated May 2011. The Health and Safety Plan (HASP) presents the specific health and safety protocols associated with the activities planned for the BCP Site.

# 7.0 CITIZEN PARTICIPATION

The Citizen Participation (CP) component for the Former Signore Facility BCP Site is discussed in the "Brownfield Cleanup Program, Citizen Participation Plans, Former Signore Facility, 55 Jefferson Street, Village of Ellicottville, Cattaraugus County, New York, Site Number: C905034" dated March 2011. The CP Plan outlines how members of the affected and interested public are provided with information about how NYSDEC will inform and involve them during the investigation and remediation of the Site. Information such as project contacts, document repositories, site contact lists, and CP activities are provided in the CP Plan.

#### 8.0 SCHEDULE

The following schedule is proposed for the IRM field activities and IRM Report preparation.

Activity:Anticipated Date:Perform Soil & EDC Injection Activities:August 2013Pilot Test Groundwater Sampling Event 1Fall 2013Resubmittal of SRI/AA Report with IRM ActivitiesFall 2013 (after 1st event)Pilot Test Groundwater Sampling Event 2Late Spring 2014



# **TABLES**

# Table 1 Estimated IRM Analytical Testing Program Summary Former Signore Facility BCP Site No. C905034 Ellicottville, New York

|                      |             |             |                            | V                    |
|----------------------|-------------|-------------|----------------------------|----------------------|
| Location             | VOCs<br>TCL | VOCs<br>TCL | Waste<br>Characterization  | EDC<br>MONITORING ** |
| AOC-2                | SOIL        | WATER       | Soil                       | Water                |
| Confirmatory         | 5           | 1           | 1                          | 0                    |
| Duplicate            | 1           | 0           | 0                          | 0                    |
| MS/MSD               | 0           | 0           | 0                          | 0                    |
| Rinsate              | 0           | 0           | 0                          | 0                    |
| Total                | 6           | 1           | 1                          | 0                    |
| AOC-3                |             |             |                            | Villay said          |
| Confirmatory         | 14          | 1           | 4                          | 0                    |
| Duplicate            | 1           | 0           | 0                          | 0                    |
| MS/MSD               | 2           | 0           | 0                          | 0                    |
| Rinsate              | 0           | 0 -         | 0                          | 0                    |
| Total                | 17          | 0           | 4                          | 0                    |
| EDC IINJECTION MONIT | TORING      |             | Minimum at the fill of the |                      |
| Various <sup>1</sup> | 0           | 12          | 0                          | 12                   |
| Duplicate            | 0           | 0           | 0                          | 0                    |
| MS/MSD               | 0           | 0           | 0                          | 0                    |
| Rinsate              | 0           | 0           | 0                          | 0                    |
| Total                | 0           | 12          | 0                          | 12                   |
|                      | 23          | 13          | 5                          | 12                   |

# Notes:

MS/MSD - Matrix Spike/Matrix Spike Duplicate.

TCL VOCs - Target Compound List Volatile Organic Compounds.

TCL SVOCs - Target Compound List Semi-volatile Organic Compounds.

TAL Metals - Target Analyte List Metals.

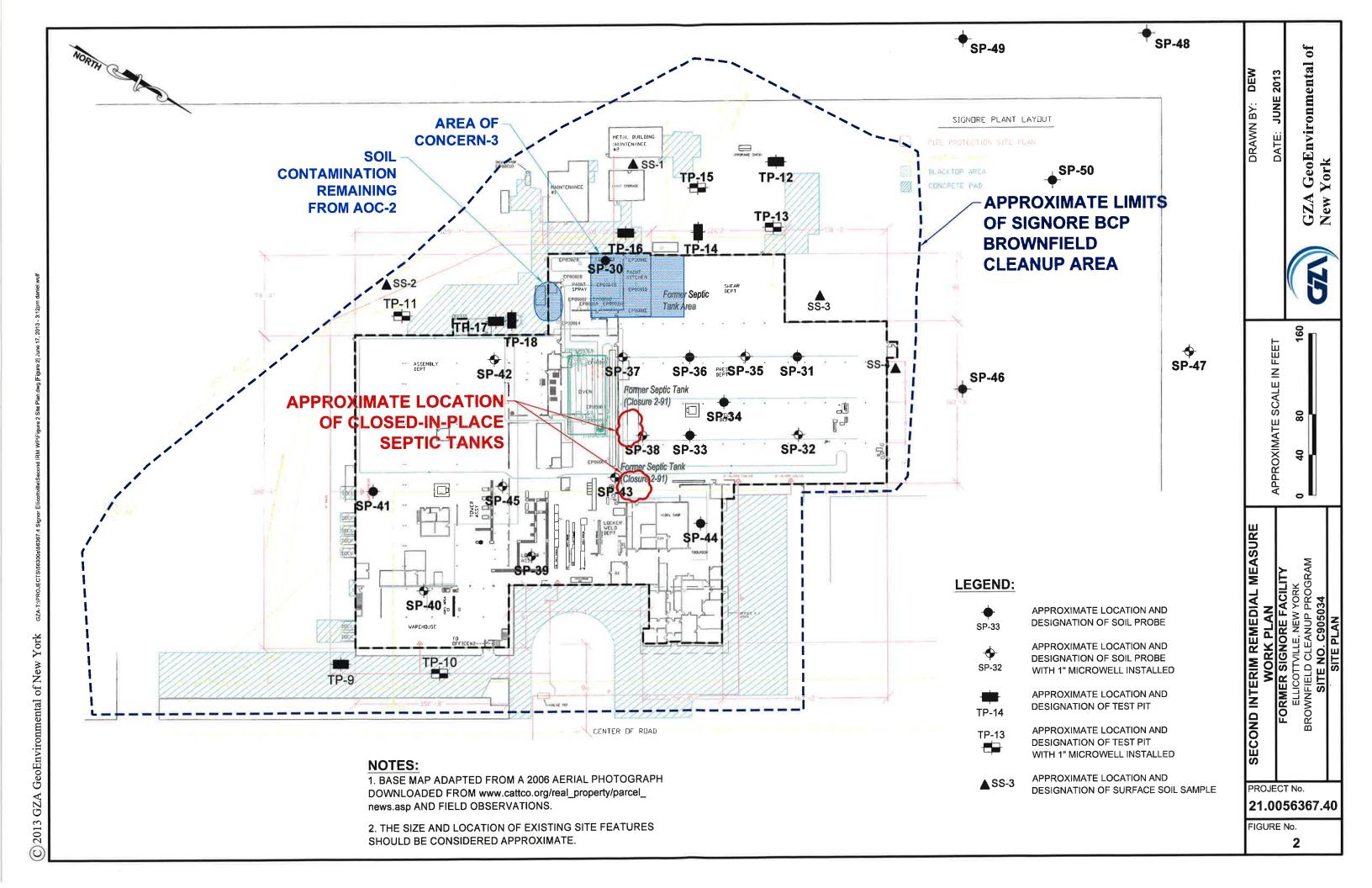
TCL PCBs - Target Compound List Polychlorinated Biphenyls.

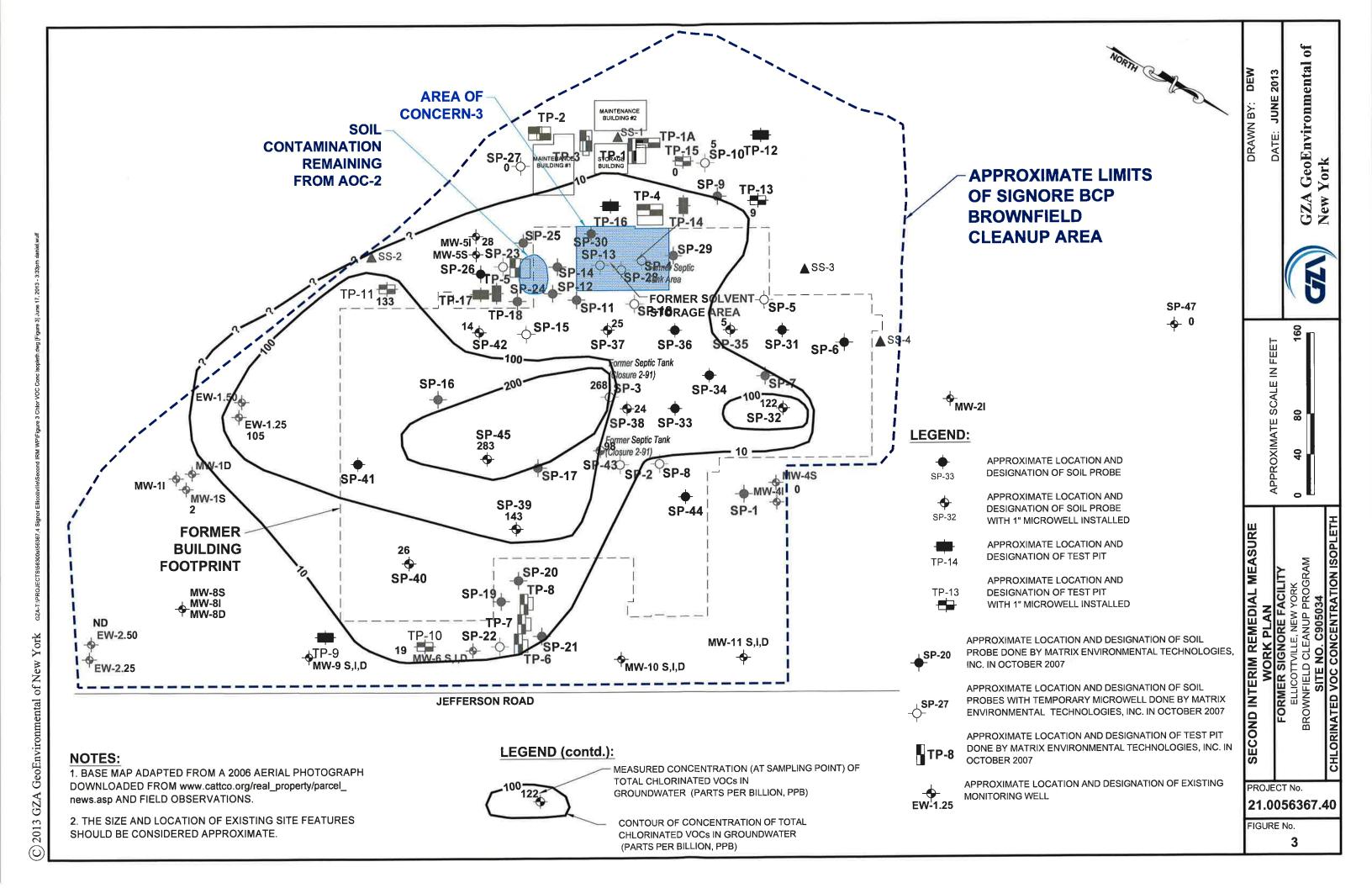
\* = Waste Characterization will be based on disposal facility requirements

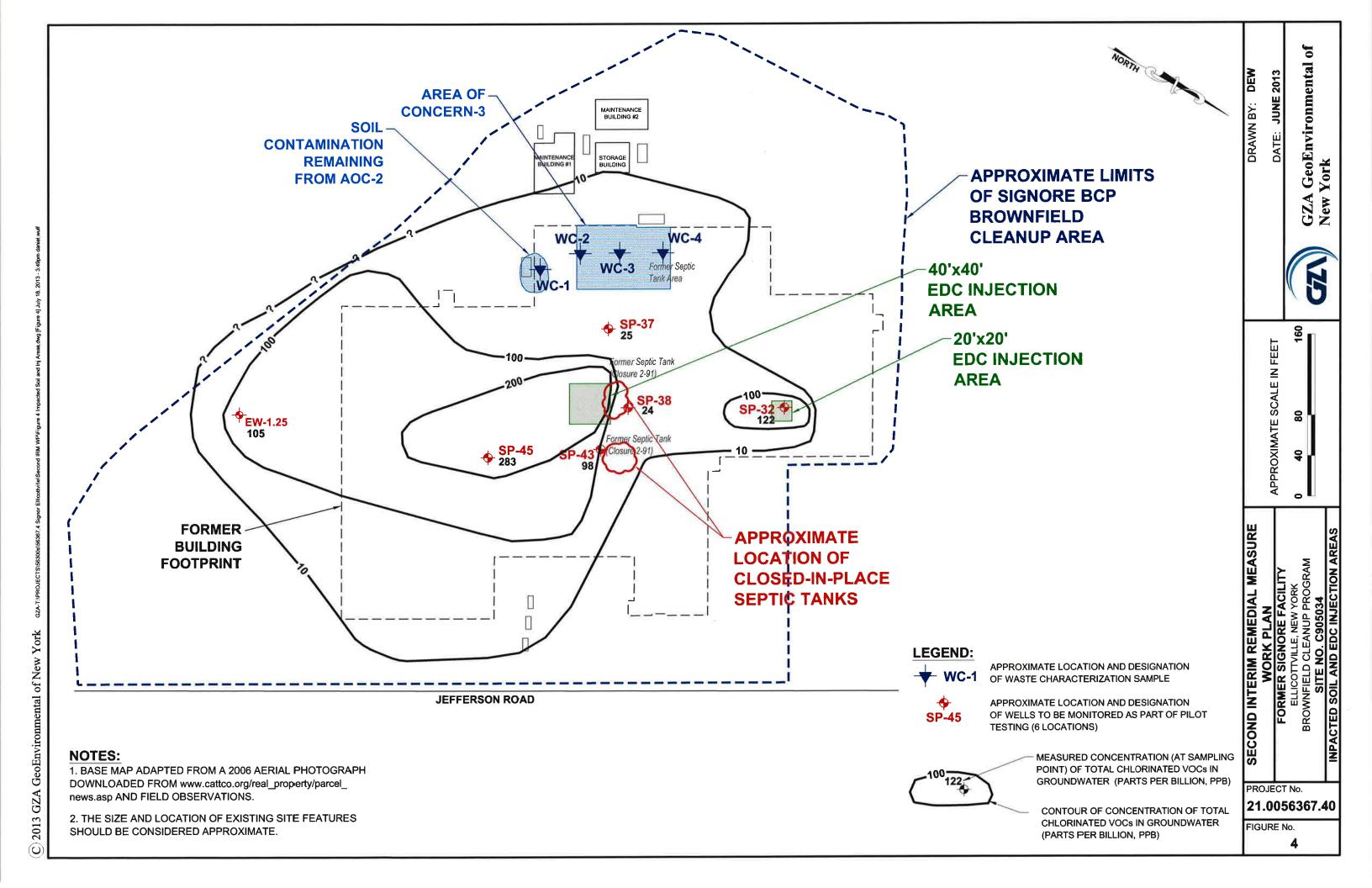
\*\* = Parameters will include list provided in Section 2.4.



# **FIGURES**









# APPENDIX A

# Table 3

# Soil Analytical Testing Results Summary Signore Facility

55-57 Jefferson Ellicottville, New York

| _                         | Unrestricted Use  | Restricted S    | Soil Cleanup Object | ives (SCO) | SP - 3 14-16ft. | SP - 4 10-12ft. | SP - 2 10-12ft. | SP - 1 18-20ft. | TP-1 9-11ft. | TP - 2 6-7ft. | TP - 4 9ft. | TP - 1A 9ft. | TP - 5 7ft. | TP - 5 9.5ft. | TP - 4 10-11ft. | TP - 5 12ft. | TP - 6 7-8ft. | TP - 7 8ft. | TP - 7 7-8ft. |
|---------------------------|-------------------|-----------------|---------------------|------------|-----------------|-----------------|-----------------|-----------------|--------------|---------------|-------------|--------------|-------------|---------------|-----------------|--------------|---------------|-------------|---------------|
| Parameter                 | Soil Cleanup      | Restricted      | Restricted          | Restricted | 10/02/2007      | 10/02/2007      | 10/02/2007      | 10/02/2007      | 10/03/2007   | 10/03/2007    | 10/03/2007  | 10/03/2007   | 10/03/2007  | 10/03/2007    | 10/03/2007      | 10/03/2007   | 10/03/2007    | 10/03/2007  | 10/03/2007    |
|                           | Objectives        | Residential     | Commercial          | Industrial | Result          | Result          | Result          | Result          | Result       | Result        | Result      | Result       | Result      | Result        | Result          | Result       | Result        | Result      | Result        |
| Volatile Organic Compour  | ds - EPA Method 8 | 260 TCL (ug/kg) |                     |            |                 |                 |                 |                 |              |               |             |              |             |               |                 |              |               |             |               |
| 1,1-Dichloroethene        | 330               | 100,000         | 500,000             | 1,000,000  | <               | <               | <               | <               | <            | 140           | <           | <            | <           | <             | <               | <            | <             | <           | <             |
| 1,1-Dichloroethane        | 270               | 26,000          | 240,000             | 480,000    | <               | <               | <               | <               | <            | 260           | <           | <            | <           | <             | <               | <            | <             | <           | <             |
| 1,1,1-Trichloroethane     | 680               | 100,000         | 500,000             | 1,000,000  | <               | <               | <               | <               | <            | 520           | <           | <            | <           | <             | <               | <            | <             | <           | <             |
| Benzene                   | 60                | 4,800           | 44,000              | 89,000     | <               | <               | <               | <               | <            | <             | <           | <            | <           | <             | <               | <            | <             | 800         | 2,900         |
| Trichloroethene           | 470               | 21,000          | 200,000             | 400,000    | 150             | <               | 73              | 130             | <            | <             | <           | <            | <           | <             | <               | <            | <             | <           | <             |
| Toluene                   | 700               | 100,000         | 500,000             | 1,000,000  | <               | <               | <               | <               | <            | <             | <           | <            | 330,000     | 250,000       | <               | 13,000       | <             | 390         | 29,000        |
| Tetrachloroethene         | 1,300             | 19,000          | 150,000             | 300,000    | 100             | <               | 220             | 64              | <            | <             | <           | <            | <           | <             | <               | <            | <             | <           | <             |
| Ethylbenzene              | 1,000             | 41,000          | 390,000             | 780,000    | <               | 78,000          | <               | <               | <            | <             | <           | <            | 38,000      | 32,000        | <               | 1,900        | <             | 4,300       | 16,000        |
| m&p-Xylene                | 260               | 100,000         | 500,000             | 1,000,000  | <               | 310,000         | <               | <               | <            | <             | <           | <            | 160,000     | 160,000       | <               | 9,900        | 69            | 22,000      | 81,000        |
| o-Xylene                  | 260               | 100,000         | 500,000             | 1,000,000  | <               | 130,000         | <               | <               | <            | <             | <           | <            | 49,000      | 56,000        | <               | 1,800        | <             | 4,700       | 27,000        |
| sopropylbenzene           | NV                | NV              | NV                  | NV         | <               | 34,000          | <               | <               | <            | <             | <           | <            | 1,500       | 1,800         | <               | <            | <             | 380         | 1,200         |
| n-Propylbenzene           | 3,900             | 100,000         | 500,000             | 1,000,000  | <               | 250,000         | <               | <               | <            | <             | <           | <            | 1,100       | 1,300         | <               | <            | 65            | 2,200       | 6,400         |
| 1,3,5-Trimethylbenzene    | 8,400             | 52,000          | 190,000             | 380,000    | <               | 550,000         | <               | <               | <            | <             | <           | <            | 630         | 930           | <               | <            | 70            | 6,100       | 20,000        |
| 1,2,4-Trimethylbenzene    | 3,600             | 52,000          | 190,000             | 380,000    | <               | 1,400,000       | <               | <               | <            | <             | <           | <            | 1,000       | 1,400         | <               | <            | 140           | 19,000      | 53,000        |
| sec-Butylbenzene          | 11,000            | 100,000         | 500,000             | 1,000,000  | <               | 21,000          | <               | <               | <            | <             | <           | <            | <           | <             | <               | <            | <             | 240         | 780           |
| o-Isopropyltoluene        | NV                | NV              | NV                  | NV         | <               | 26,000          | <               | <               | <            | <             | <           | <            | <           | <             | <               | <            | <             | 430         | 1,400         |
| n-Butylbenzene            | 12,000            | 100,000         | 500,000             | 1,000,000  | <               | 32,000          | <               | <               | <            | <             | <           | <            | <           | <             | <               | <            | <             | 950         | 2,500         |
| Naphthalene               | 12,000            | 100,000         | 500,000             | 1,000,000  | <               | <               | <               | <               | <            | 99            | <           | <            | <           | <             | <               | <            | <             | 1,200       | 3,200         |
| Total VOCs                |                   |                 |                     |            | 250             | 2,831,000       | 293             | 194             |              | 1,019         |             |              | 581,230     | 503,430       |                 | 26,600       | 344           | 62,690      | 244,380       |
| Semi-Volatile Organic Cor | npounds - EPA Met | hod 8270 STAR   | S (ug/kg)           |            |                 |                 |                 |                 |              |               |             |              |             |               |                 |              |               |             |               |
| Naphthalene               | 12,000            | 100,000         | 500,000             | 1,000,000  | NT              | NT              | NT              | NT              | NT           | NT            | NT          | NT           | NT          | NT            | NT              | NT           | <             | 2,000       | 730           |
| 2-Methylnaphthalene       | NV                | NV              | NV                  | NV         | NT              | NT              | NT              | NT              | NT           | NT            | NT          | NT           | NT          | NT            | NT              | NT           | <             | 4,000       | 1,300         |
| Phenanthrene              | 100,000           | 100,000         | 500,000             | 1,000,000  | NT              | NT              | NT              | NT              | NT           | NT            | NT          | NT           | NT          | NT            | NT              | NT           | <             | 450         | <             |

|                           | Unrestricted Use   | Restricted S    | Soil Cleanup Objecti | ves (SCO)  | TP - 8 7-8ft. | SP - 16 10-12 | SP - 9 4-6ft. | SP - 13 10-12 | SP - 15 14-16 | SP-19 2-4  | SP-20 8-10 | SP-21 8-10 | SP-22 8-10 | SP-23 8-10 | SP-24 8-10 | SP-25 8-10 | SP-26 8-10 | SP-28 8-10 | SP-29 8-10 |
|---------------------------|--------------------|-----------------|----------------------|------------|---------------|---------------|---------------|---------------|---------------|------------|------------|------------|------------|------------|------------|------------|------------|------------|------------|
| Parameter                 | Soil Cleanup       | Restricted      | Restricted           | Restricted | 10/03/2007    | 10/03/2007    | 10/03/2007    | 10/03/2007    | 10/03/2007    | 10/05/2007 | 10/05/2007 | 10/05/2007 | 10/05/2007 | 10/05/2007 | 10/05/2007 | 10/05/2007 | 10/05/2007 | 10/05/2007 | 10/05/2007 |
|                           | Objectives         | Residential     | Commercial           | Industrial | Result        | Result        | Result        | Result        | Result        | Result     | Result     | Result     | Result     | Result     | Result     | Result     | Result     | Result     | Result     |
| Volatile Organic Compound | ls - EPA Method 82 | 260 TCL (ug/kg) |                      |            |               |               |               |               |               |            |            |            |            |            |            |            |            |            |            |
| 1,1-Dichloroethene        | 330                | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | <             | <             | <          | <          | <          | <          | <          | <          | <          | <          | <          | <          |
| 1,1-Dichloroethane        | 270                | 26,000          | 240,000              | 480,000    | <             | <             | <             | <             | <             | <          | <          | <          | <          | <          | <          | <          | <          | <          | <          |
| 1,1,1-Trichloroethane     | 680                | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | <             | <             | <          | <          | <          | <          | <          | <          | <          | <          | <          | <          |
| Benzene                   | 60                 | 4,800           | 44,000               | 89,000     | <             | <             | <             | <             | <             | <          | <          | <          | <          | <          | <          | <          | <          | <          | <          |
| Trichloroethene           | 470                | 21,000          | 200,000              | 400,000    | <             | 260           | <             | <             | <             | <          | <          | <          | <          | <          | <          | <          | <          | <          | 60         |
| Toluene                   | 700                | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | <             | <             | <          | 86         | <          | <          | <          | <          | <          | <          | <          | <          |
| Tetrachloroethene         | 1,300              | 19,000          | 150,000              | 300,000    | <             | 1,200         | <             | <             | <             | <b>~</b>   | <          | <          | <b>~</b>   | <          | <          | <          | <          | <          | <          |
| Ethylbenzene              | 1,000              | 41,000          | 390,000              | 780,000    | <             | <             | <             | 2,300         | 19,000        | <          | 1,000      | <          | <          | <          | <          | <          | <          | 34,000     | <          |
| m&p-Xylene                | 260                | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | 8,700         | 33,000        | <          | 3,600      | <          | <          | <          | <          | <          | <          | 140,000    | <          |
| o-Xylene                  | 260                | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | 2,300         | 900           | <          | 210        | <          | <          | <          | <          | <          | <          | 44,000     | <          |
| Isopropylbenzene          | NV                 | NV              | NV                   | NV         | <             | <             | <             | 2,000         | 940           | <          | 660        | <          | <          | <          | <          | <          | <          | 21,000     | <          |
| n-Propylbenzene           | 3,900              | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | 17,000        | 570           | <          | 730        | <          | <          | <          | <          | <          | <          | 150,000    | <          |
| 1,3,5-Trimethylbenzene    | 8,400              | 52,000          | 190,000              | 380,000    | <             | <             | <             | 34,000        | <             | <          | 1,700      | <          | <          | <          | <          | <          | <          | 350,000    | <          |
| 1,2,4-Trimethylbenzene    | 3,600              | 52,000          | 190,000              | 380,000    | <             | <             | <             | 90,000        | <             | <b>~</b>   | 4,700      | <          | <b>~</b>   | <          | <          | <          | <          | 910,000    | 96         |
| sec-Butylbenzene          | 11,000             | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | 1,600         | <             | <          | 110        | <          | <          | <          | <          | <          | <          | 13,000     | <          |
| p-lsopropyltoluene        | NV                 | NV              | NV                   | NV         | <             | <             | <             | 2,300         | <             | <          | 220        | <          | <          | <          | <          | <          | <          | 20,000     | <          |
| n-Butylbenzene            | 12,000             | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | 2,300         | <             | <          | 520        | <          | <          | <          | <          | <          | <          | 19,000     | <          |
| Naphthalene               | 12,000             | 100,000         | 500,000              | 1,000,000  | <             | <             | <             | <             | <             | <          | 1,100      | <          | <          | <          | <          | <          | <          | <          | <          |
| Total VOCs                |                    |                 |                      |            | 0             | 1,460         | 0             | 162,500       | 54,410        |            | 14,636     |            |            |            |            |            |            | 1,701,000  | 156        |
| Semi-Volatile Organic Com | pounds - EPA Meth  | nod 8270 STAR   | S (ug/kg)            |            |               |               |               |               |               |            |            |            |            |            |            |            |            |            |            |
| Naphthalene               | 12,000             | 100,000         | 500,000              | 1,000,000  | <             | NT            | NT            | NT            | NT            | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         |
| 2-Methylnaphthalene       | NV                 | NV              | NV                   | NV         | <             | NT            | NT            | NT            | NT            | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         |
| Phenanthrene              | 100,000            | 100,000         | 500,000              | 1,000,000  | <             | NT            | NT            | NT            | NT            | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         | NT         |

- Compounds detected in one or more samples are presented on this table. Refer to Attachment C for list of all compounds included in analysis.
   Analytical testing completed by GZA GeoEnvironmental Laboratory.
   ug/kg = part per billion and mg/kg = parts per million.
   cindicates compound was not detected.

- 5. Bold indicates value exceeds the Unrestricted Use Soil Cleanup Objectives
- Blue shading indicates value exceeds the Restricted Residential Use Soil Cleanup Objectives
   Yellow shading indicates value exceeds the Restricted Commercial Use Soil Cleanup Objectives
   Red shading indicates value exceeds the Restricted Industrial Use Soil Cleanup Objectives



# APPENDIX B



| ✓ Final Repo | ort    |
|--------------|--------|
| Re-Issued    | Report |
| Revised R    | eport  |

# Laboratory Report

GZA GeoEnvironmental of NY Buffalo 535 Washington Street, 11th Floor

Buffalo, NY 14203

Attn: Chris Boron

Work Order: M1089

Project: Former Signore Facility

Project #:

| Laboratory ID | Client Sample ID | <u>Matrix</u> | Date Sampled    | Date Received   |
|---------------|------------------|---------------|-----------------|-----------------|
| M1089-01      | WC-2-061313      | Soil          | 13-Jun-13 15:00 | 03-Jul-13 12:17 |
| M1089-02      | WC-3-061313      | Soil          | 13-Jun-13 15:00 | 03-Jul-13 12:17 |

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. The results relate only to the samples(s) as received. This report may not be reproduced, except in full, without written approval from Spectrum Analytical.

All applicable NELAC or USEPA CLP requirments have been meet.

Spectrum Analytical (Rhode Island) is accredited under the National Environmental Laboratory Approval Program (NELAP) and DoD Environmental Laboratory Accreditation Program (ELAP), holds Organic and Inorganic contracts under the USEPA CLP Program and is certified under several states. The current list of our laboratory approvals and certifications is available on the Certifications page on our web site at www.spectrum-analytical.com.

Please contact the Laboratory or Technical Director at 401-732-3400 with any questions regarding the data contained in the laboratory report.

Department of Defense N/A PH-0153 Connecticut Delaware N/A Florida E87664 Maine 2007037 Massachusetts M-RI907 New Hampshire 2631 New Jersey RI001 New York 11522 North Carolina 581 Rhode Island LAI00301 P330-08-00023 USDA USEPA - ISM EP-W-09-039 USEPA - SOM EP-W-11-033





Authorized by:

Yihai Ding Laboratory Director



\* Data Summary Pack \*

# **New York State Department of Environmental Conservation Sample Identification and Analytical Requirements Summary**

Project Name : Former Signore Facility

SDG: M1089

|                       |                         |                   | Analy              | tical Requirement | ts       |       |
|-----------------------|-------------------------|-------------------|--------------------|-------------------|----------|-------|
| Customer<br>Sample ID | Laboratory<br>Sample ID | MSVOA<br>Method # | MSSEMI<br>Method # | GC*<br>Method #   | ME       | Other |
| WC-2-061313           | M1089-01                |                   | SW8270_S           |                   | SW6010_S |       |
| WC-2-061313           | M1089-01                |                   |                    |                   | SW7471   |       |
| WC-3-061313           | M1089-02                |                   | SW8270_S           |                   | SW6010_S |       |
| WC-3-061313           | M1089-02                |                   |                    |                   | SW7471   |       |

**Page 1** 07/12/2013 13:29

# New York State Department of Environmental Conservation Sample Preparation and Analysis Summary MSSEMI

Project Name: Former Signore Facility SDG: M1089

| Laboratory<br>Sample ID | Matrix | Date<br>Collected | Date Received<br>By Lab | Date<br>Extracted | Date<br>Analyzed |
|-------------------------|--------|-------------------|-------------------------|-------------------|------------------|
| SW8270_S                |        |                   |                         |                   |                  |
| M1089-01A               | SL     | 6/13/2013         | 7/3/2013                | 7/5/2013          | 7/5/2013         |
| M1089-02A               | SL     | 6/13/2013         | 7/3/2013                | 7/5/2013          | 7/5/2013         |

Page 3 07/12/2013 13:29

# New York State Department of Environmental Conservation Sample Preparation and Analysis Summary MSSEMI

Project Name: Former Signore Facility SDG: M1089

| Laboratory<br>Sample ID | Matrix | Analytical<br>Protocol | Extraction<br>Method | Auxiliary<br>Cleanup | Dil/Conc<br>Factor |
|-------------------------|--------|------------------------|----------------------|----------------------|--------------------|
| SW8270_S                |        |                        |                      |                      |                    |
| M1089-01A               | SL     | SW8270_S               | 3550B                | NA                   | 1                  |
| M1089-02A               | SL     | SW8270_S               | 3550B                | NA                   | 1                  |

**Page 6** 07/12/2013 13:29

# New York State Department of Environmental Conservation Sample Preparation and Analysis Summary ME

Project Name : Former Signore Facility

SDG: M1089

| Laboratory |        | Metals    | Date Received | Date     |
|------------|--------|-----------|---------------|----------|
| Sample ID  | Matrix | Requested | By Lab        | Analyzed |
| SW6010_S   |        |           |               |          |
| M1089-01A  | SL     | SW6010_S  | 7/3/2013      | 7/9/2013 |
| M1089-02A  | SL     | SW6010_S  | 7/3/2013      | 7/9/2013 |
| SW7471     |        |           |               |          |
| M1089-01A  | SL     | SW7471    | 7/3/2013      | 7/9/2013 |
| M1089-02A  | SL     | SW7471    | 7/3/2013      | 7/9/2013 |

**Page 8** 07/12/2013 13:29

# ge EQUIIS\_4\_NYSDEC

Sectrum Analytical Inc. - North Kingstown RI -- Rhode Island Division Client ID: GZA\_BUFFALO

**Project:** Former Signore Facility

Report Level: ASP-B Special Program: HC Due: 07/12/13 Fax Due: 07/09/13 Case: SDG:

EDD: CLF Fax Report:

WorkOrder: M1089

Comments: CC pdf and EDD to John Beninati (john.beninati@gza.com). No charge for Trip Blank. FORMERLY M0976-REANLYSIS REQUESTED BY CLIENT **PO:** 21.0056491.00 WO Name: Former Signore Facility Location: GZA\_SINGNORE,

| Lab Samp ID | Lab Samp ID Client Sample ID | Collection Date             | Date Recv'd | Matrix | Matrix Test Code | Samp / Lab Test Comments                | HF HT MS SEL Storage | EL Sto  | rage |
|-------------|------------------------------|-----------------------------|-------------|--------|------------------|---|----------------------|---------|------|
| M1089-01A   | WC-2-061313                  | 06/13/2013 15:00 07/03/2013 | 07/03/2013  | Soil   | PMoist           | USE PMOIST FROM M0976-02A /             |                      | A1      |      |
| M1089-01A   | WC-2-061313                  | 06/13/2013 15:00 07/03/2013 | 07/03/2013  | Soil   | SW6010_S         | USE PMOIST FROM M0976-02A / TAL         | >                    | ,<br>A1 |      |
| M1089-01A   | WC-2-061313                  | 06/13/2013 15:00 07/03/2013 | 07/03/2013  | Soil   | SW7471           | USE PMOIST FROM M0976-02A / TAL         |                      | A1      |      |
| M1089-01A   | WC-2-061313                  | 06/13/2013 15:00            | 07/03/2013  | Soil   | SW8270_S         | USE PMOIST FROM M0976-02A /<br>8270_BN, | >                    | ,<br>A1 |      |
| M1089-02A   | WC-3-061313                  | 06/13/2013 15:00 07/03/2013 | 07/03/2013  | Soil   | PMoist           | USE PMOIST FROM M0976-03A /             |                      | A1      |      |
| M1089-02A   | WC-3-061313                  | 06/13/2013 15:00 07/03/2013 | 07/03/2013  | Soil   | SW6010_S         | USE PMOIST FROM M0976-03A / TAL         | >                    | A1      |      |
| M1089-02A   | WC-3-061313                  | 06/13/2013 15:00            | 07/03/2013  | Soil   | SW7471           | USE PMOIST FROM M0976-03A / TAL         |                      | A1      |      |
| M1089-02A   | WC-3-061313                  | 06/13/2013 15:00 07/03/2013 | 07/03/2013  | Soil   | SW8270_S         | USE PMOIST FROM M0976-03A /<br>8270_BN, | <b>&gt;</b>          | ,<br>A1 |      |

Lab Client Rep: Veronica E Brizard

Page 01 of 01

HT = Test logged in but has been placed on hold



\* Semivolatile Organics \*

# REPORT NARRATIVE

Spectrum Analytical, Inc. Featuring Hanibal Technology, RI Division.

Client: GZA GeoEnvironmental of NY Buffalo

**Project: Former Signore Facility** 

Laboratory Workorder / SDG #: M1089

**SW846 8270D, SVOA by GC-MS** 

#### I. SAMPLE RECEIPT

No exceptions or unusual conditions were encountered unless a Sample Condition Notification Form, or other record of communication is included with the Sample Receipt Documentation.

# II. HOLDING TIMES

# A. Sample Preparation:

All samples were prepared within the method-specified holding times with the following exceptions:

WC-2-061313 (M1089-01A) exceed by-8 Days WC-3-061313 (M1089-02A) exceed by-8 Days

Please note these two samples were collected on 6/13. Additional tests were added on 7/3.

# B. Sample Analysis:

All samples were analyzed within the method-specified holding times.

# III. METHODS

Samples were analyzed following procedures in laboratory test code: SW846 8270D

#### IV. PREPARATION

Soil Samples were prepared following procedures in laboratory test code: SW3550B

M1089 Page 3 of 37

#### V. INSTRUMENTATION

The following instrumentation was used

Instrument Code: S4

Instrument Type: GCMS-SEMI Description: HP6890 / HP5973 Manufacturer: Hewlett-Packard

Model: 6890N / 5973N

# VI. ANALYSIS

# A. Calibration:

Calibrations met the method/SOP acceptance criteria.

#### B. Blanks:

All method blanks were within the acceptance criteria.

# C. Surrogates:

Surrogate standard percent recoveries were within the QC limits.

# D. Spikes:

# 1. Laboratory Control Spikes (LCS):

Percent recoveries for lab control samples were within the QC limits with the following exceptions. Please note that most test procedures allow for several compounds outside of the QC limits for the LCS, although this may indicate a bias for this specific compound.

Replicate RPDs were within the advisory QC limits with the exception of the following:

(LCSD-72594), Relative Percent Difference is greater than RPD limit for 3,3´-Dichlorobenzidine, 4-Chloroaniline.

# Matrix Spike / Matrix Spike Duplicate (MS/MSD):

No client-requested MS/MSD analyses were included in this SDG.

# E. Internal Standards:

M1089 Page 4 of 37

Internal standard peak areas were within the QC limits.

### F. Dilutions:

No sample in this SDG required analysis at dilution.

## G. Samples:

No other unusual occurrences were noted during sample analysis.

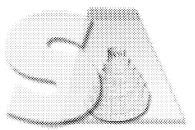
## H. Manual Integration

No manual integrations were performed on any sample or standard.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and Spectrum, both technically and for completeness, except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

| tSigned: | J-W       |  |
|----------|-----------|--|
| Nate:    | 7/12/2013 |  |

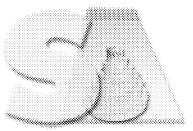
M1089 Page 5 of 37



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

# Data Flag/Qualifiers:

- U Not Detected. This compound was analyzed-for but not detected. For most analyses the reporting limit (lowest standard concentration) is the value listed. For Department of Defense programs, this is the Limit of Detection (LOD).
- J This flag indicates an estimated value due to either
  - the compound was detected below the reporting limit, or
  - estimated concentration for Tentatively Identified Compound
- B This flag indicates the compound was also detected in the associated Method Blank. The B flag has an alternative meaning for Inorganics analyses reported using CLP ILM-type metals forms, indicating a "trace" concentration below the reporting limit and equal to or above the detection limit.
- D For Organics analysis, this flag indicates the compound concentration was obtained from a secondary dilution analysis
- E This flag indicates the compound concentration exceeded the Calibration Range. The E flag has an alternative meaning for Inorganics analyses reported using CLP metals forms, indicating an estimated concentration due to the presence of interferences, as determined by the serial dilution analysis.
- P This flag is used for pesticides/PCB/herbicide compound when there is a greater than 40% difference for detected concentration between the two GC columns used for primary and confirmation analyses. This difference typically indicates an interference, causing one value to be unusually high. The **lower** of the two values is generally reported on the Form 1, and both values reported on the Form 10.
- A Used to flag semivolatile organic Tentatively Identified Compound library search results for compounds identified as aldol condensation byproducts.
- N Used to flag results for volatile and semivolatile Organics analysis Tentatively Identified Compounds where an analyte has passed the identification criteria, and is considered to be positively identified. For Inorganics analysis the N flag indicates the matrix spike recovery falls outside of the control limit.
- \* For Inorganics analysis the \* flag indicates Relative Percent Difference for duplicate analyses is outside of the control limit.



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

# **Sample ID Suffixes**

- DL Diluted analysis. The sample was diluted and reanalyzed. The DL may be followed by a digit if more than one diluted reanalysis is provided. The DL suffix is not attached to an analysis initially performed at dilution, only to reanalyses performed at dilution
- RE Reanalysis. Appended to the client sample ID to indicate a reextraction and reanalysis or a reanalysis of the original sample extract.
- RA Reanalysis. Appended to the laboratory sample ID indicates a reanalysis of the original sample extract.
- RX Reextraction. Appended to the laboratory sample ID indicates a reextraction of the sample.
- MS Matrix Spike.
- MSD Matrix Spike Duplicate
- DUP Duplicate analysis
- SD Serial Dilution
- PS Post-digestion or Post-distillation spike. For metals or inorganic analyses

# 1D - FORM I SV-1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO. WC-2-061313

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: M1089-01A      |
| Sample wt/vol:15.2 (g/mL) G               | Lab File ID: S4F4234.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: Decanted: (Y/N) N             | Date Received: 07/03/2013     |
| Concentrated Extract Volume:1000 (uL)     | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

|           |                              | CONCENTRATION UNITS:  |   |
|-----------|------------------------------|-----------------------|---|
| CAS NO.   | COMPOUND                     | (ug/L or ug/Kg) UG/KG | Q |
| 111-44-4  | Bis(2-chloroethyl)ether      | 370                   | Ū |
|           | 1,3-Dichlorobenzene          | 370                   | Ū |
| 106-46-7  | 1,4-Dichlorobenzene          | 370                   | Ū |
| 95-50-1   | 1,2-Dichlorobenzene          | 370                   | Ū |
| 108-60-1  | 2,2´-oxybis(1-Chloropropane) | 370                   | Ū |
|           | Hexachloroethane             | 370                   | U |
| 98-95-3   | Nitrobenzene                 | 370                   | U |
| 78-59-1   | Isophorone                   | 370                   | U |
| 120-82-1  | 1,2,4-Trichlorobenzene       | 370                   | U |
| 91-20-3   | Naphthalene                  | 370                   | U |
| 106-47-8  | 4-Chloroaniline              | 370                   | U |
| 111-91-1  | Bis(2-chloroethoxy)methane   | 370                   | U |
| 87-68-3   | Hexachlorobutadiene          | 370                   | U |
| 91-57-6   | 2-Methylnaphthalene          | 370                   | U |
| 77-47-4   | Hexachlorocyclopentadiene    | 370                   | U |
| 91-58-7   | 2-Chloronaphthalene          | 370                   | U |
| 88-74-4   | 2-Nitroaniline               | 750                   | U |
| 131-11-3  | Dimethylphthalate            | 370                   | U |
| 208-96-8  | Acenaphthylene               | 370                   | U |
| 606-20-2  | 2,6-Dinitrotoluene           | 370                   | U |
| 99-09-2   | 3-Nitroaniline               | 750                   | U |
| 83-32-9   | Acenaphthene                 | 370                   | U |
| 132-64-9  | Dibenzofuran                 | 370                   | U |
| 121-14-2  | 2,4-Dinitrotoluene           | 370                   | U |
| 84-66-2   | Diethylphthalate             | 370                   | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether   | 370                   | U |
| 86-73-7   | Fluorene                     | 370                   | U |
| 100-01-6  | 4-Nitroaniline               | 750                   | U |
| 101-55-3  | 4-Bromophenyl-phenylether    | 370                   | U |
| 118-74-1  | Hexachlorobenzene            | 370                   | U |
| 85-01-8   | Phenanthrene                 | 370                   | U |
| 120-12-7  | Anthracene                   | 370                   | U |
| 86-74-8   | Carbazole                    | 370                   | U |
| 206-44-0  | Fluoranthene                 | 370                   | U |
| 129-00-0  |                              | 370                   | U |
| 85-68-7   | Butylbenzylphthalate         | 370                   | U |

# 1E - FORM I SV-2 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO. WC-2-061313

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: M1089-01A      |
| Sample wt/vol:15.2 (g/mL) G               | Lab File ID: S4F4234.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: Decanted: (Y/N) N             | Date Received: 07/03/2013     |
| Concentrated Extract Volume:1000 (uL)     | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

| CAS NO.  | COMPOUND                   | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG | Q |
|----------|----------------------------|--|---|
| 91-94-1  | 3,3´-Dichlorobenzidine     | 370  | U |
| 56-55-3  | Benzo(a)anthracene         | 370  | U |
| 218-01-9 | Chrysene                   | 370  | U |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 370  | U |
| 205-99-2 | Benzo(b)fluoranthene       | 370  | U |
|          | Benzo(k)fluoranthene       | 370  | U |
| 50-32-8  | Benzo(a)pyrene             | 370  | U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene     | 370  | U |
|          | Dibenzo(a,h)anthracene     | 370  | U |
| 191-24-2 | Benzo(g,h,i)perylene       | 370  | U |

# 1D - FORM I SV-1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

WC-3-061313

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: M1089-02A      |
| Sample wt/vol:15.3 (g/mL) G               | Lab File ID: S4F4235.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: 8.9 Decanted: (Y/N) N         | Date Received: 07/03/2013     |
| Concentrated Extract Volume: 1000 (uL)    | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

|           |                              | CONCENTRATION UNITS:  |   |
|-----------|------------------------------|-----------------------|---|
| CAS NO.   | COMPOUND                     | (ug/L or ug/Kg) UG/KG | Q |
| 111-44-4  | Bis(2-chloroethyl)ether      | 360                   | U |
|           | 1,3-Dichlorobenzene          | 360                   | U |
| 106-46-7  | 1,4-Dichlorobenzene          | 360                   | U |
| 95-50-1   | 1,2-Dichlorobenzene          | 360                   | U |
| 108-60-1  | 2,2'-oxybis(1-Chloropropane) | 360                   | U |
|           | Hexachloroethane             | 360                   | U |
| 98-95-3   | Nitrobenzene                 | 360                   | U |
| 78-59-1   | Isophorone                   | 360                   | U |
| 120-82-1  | 1,2,4-Trichlorobenzene       | 360                   | U |
| 91-20-3   | Naphthalene                  | 360                   | U |
| 106-47-8  | 4-Chloroaniline              | 360                   | U |
| 111-91-1  | Bis(2-chloroethoxy)methane   | 360                   | U |
| 87-68-3   | Hexachlorobutadiene          | 360                   | U |
| 91-57-6   | 2-Methylnaphthalene          | 360                   | U |
| 77-47-4   | Hexachlorocyclopentadiene    | 360                   | U |
| 91-58-7   | 2-Chloronaphthalene          | 360                   | U |
| 88-74-4   | 2-Nitroaniline               | 720                   | U |
| 131-11-3  | Dimethylphthalate            | 360                   | U |
| 208-96-8  | Acenaphthylene               | 360                   | U |
| 606-20-2  | 2,6-Dinitrotoluene           | 360                   | U |
| 99-09-2   | 3-Nitroaniline               | 720                   | U |
| 83-32-9   | Acenaphthene                 | 360                   | U |
| 132-64-9  | Dibenzofuran                 | 360                   | U |
| 121-14-2  | 2,4-Dinitrotoluene           | 360                   | U |
| 84-66-2   | Diethylphthalate             | 360                   | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether   | 360                   | U |
| 86-73-7   | Fluorene                     | 360                   | U |
| 100-01-6  | 4-Nitroaniline               | 720                   | U |
| 101-55-3  | 4-Bromophenyl-phenylether    | 360                   | U |
| 118-74-1  | Hexachlorobenzene            | 360                   | U |
| 85-01-8   | Phenanthrene                 | 360                   | U |
| 120-12-7  | Anthracene                   | 360                   | U |
| 86-74-8   | Carbazole                    | 360                   | U |
| 206-44-0  | Fluoranthene                 | 360                   | U |
| 129-00-0  |                              | 360                   | U |
| 85-68-7   | Butylbenzylphthalate         | 360                   | U |

# 1E - FORM I SV-2 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO. WC-3-061313

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: M1089-02A      |
| Sample wt/vol:15.3 (g/mL) G               | Lab File ID: S4F4235.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: 8.9 Decanted: (Y/N) N         | Date Received: 07/03/2013     |
| Concentrated Extract Volume: 1000 (uL)    | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

| CAS NO.  | COMPOUND                   | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG | Q |
|----------|----------------------------|--|---|
| 91-94-1  | 3,3´-Dichlorobenzidine     | 360  | U |
| 56-55-3  | Benzo(a)anthracene         | 360  | U |
| 218-01-9 | Chrysene                   | 360  | U |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 360  | U |
| 205-99-2 | Benzo(b)fluoranthene       | 360  | U |
| 207-08-9 | Benzo(k)fluoranthene       | 360  | U |
| 50-32-8  | Benzo(a)pyrene             | 360  | U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene     | 360  | U |
| 53-70-3  | Dibenzo(a,h)anthracene     | 360  | U |
| 191-24-2 | Benzo(g,h,i)perylene       | 360  | U |

# 1D - FORM I SV-1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| CLIENT | SAMPLE | NO. |
|--------|--------|-----|
| MB-725 | 94     |     |
|        |        |     |
|        |        |     |

| Lab Name: SPECTRUM ANALYTICAL, INC.         | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089            | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL               | Lab Sample ID: MB-72594       |
| Sample wt/vol:15.0 (g/mL) G                 | Lab File ID: S4F4227.D        |
| Level: (LOW/MED) LOW                        | Extraction: (Type) SONC       |
| % Moisture: Decanted: (Y/N)                 | Date Received:                |
| Concentrated Extract Volume:1000 (uL)       | Date Extracted: 07/05/2013    |
| Injection Volume: 1.0 (uL) GPC Factor: 1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                     | Dilution Factor: 1.0          |

|           |                              | CONCENTRATION UNITS:  |   |
|-----------|------------------------------|-----------------------|---|
| CAS NO.   | COMPOUND                     | (ug/L or ug/Kg) UG/KG | Q |
| 111-44-4  | Bis(2-chloroethyl)ether      | 330                   | U |
|           | 1,3-Dichlorobenzene          | 330                   | U |
| 106-46-7  | 1,4-Dichlorobenzene          | 330                   | U |
| 95-50-1   | 1,2-Dichlorobenzene          | 330                   | U |
| 108-60-1  | 2,2'-oxybis(1-Chloropropane) | 330                   | U |
|           | Hexachloroethane             | 330                   | U |
| 98-95-3   | Nitrobenzene                 | 330                   | U |
| 78-59-1   | Isophorone                   | 330                   | U |
| 120-82-1  | 1,2,4-Trichlorobenzene       | 330                   | U |
| 91-20-3   | Naphthalene                  | 330                   | U |
| 106-47-8  | 4-Chloroaniline              | 330                   | U |
| 111-91-1  | Bis(2-chloroethoxy)methane   | 330                   | U |
| 87-68-3   | Hexachlorobutadiene          | 330                   | U |
| 91-57-6   | 2-Methylnaphthalene          | 330                   | U |
| 77-47-4   | Hexachlorocyclopentadiene    | 330                   | U |
|           | 2-Chloronaphthalene          | 330                   | U |
| 88-74-4   | 2-Nitroaniline               | 670                   | U |
| 131-11-3  | Dimethylphthalate            | 330                   | U |
|           | Acenaphthylene               | 330                   | U |
|           | 2,6-Dinitrotoluene           | 330                   | U |
| 99-09-2   | 3-Nitroaniline               | 670                   | U |
| 83-32-9   | Acenaphthene                 | 330                   | U |
| 132-64-9  | Dibenzofuran                 | 330                   | U |
| 121-14-2  | 2,4-Dinitrotoluene           | 330                   | U |
| 84-66-2   | Diethylphthalate             | 330                   | U |
| 7005-72-3 | 4-Chlorophenyl-phenylether   | 330                   | U |
|           | Fluorene                     | 330                   | U |
| 100-01-6  | 4-Nitroaniline               | 670                   | U |
|           | 4-Bromophenyl-phenylether    | 330                   | U |
| 118-74-1  | Hexachlorobenzene            | 330                   | U |
| 85-01-8   | Phenanthrene                 | 330                   | U |
|           | Anthracene                   | 330                   | U |
| 86-74-8   | Carbazole                    | 330                   | U |
| 206-44-0  | Fluoranthene                 | 330                   | U |
| 129-00-0  | 2                            | 330                   | U |
| 85-68-7   | Butylbenzylphthalate         | 330                   | U |

# 1E - FORM I SV-2 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| CLIENT | SAMPLE | NO. |
|--------|--------|-----|
| MB-725 | 94     |     |
|        |        |     |
|        |        |     |

| Lab Name: SPECTRUM  | A ANALYTICAL, INC.        | Contract:         |                 |
|---------------------|---------------------------|-------------------|-----------------|
| Lab Code: MITKEM    | Case No.: M1089           | Mod. Ref No.:     | SDG No.: SM1089 |
| Matrix: (SOIL/SED/W | WATER) SOIL               | Lab Sample ID:    | MB-72594        |
| Sample wt/vol:      | 15.0 (g/mL) <u>G</u>      | Lab File ID:      | S4F4227.D       |
| Level: (LOW/MED) I  | LOW                       | Extraction: (Type | e) SONC         |
| % Moisture:         | Decanted: (Y/N)           | Date Received:    |                 |
| Concentrated Extra  | ct Volume:1000 (uL)       | Date Extracted:   | 07/05/2013      |
| Injection Volume:   | 1.0 (uL) GPC Factor: 1.00 | Date Analyzed:    | 07/05/2013      |
| GPC Cleanup:(Y/N)   | И рн:                     | Dilution Factor:  | 1.0             |

| CAS NO.  | COMPOUND                   | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG | Q |
|----------|----------------------------|--|---|
| 91-94-1  | 3,3´-Dichlorobenzidine     | 330  | U |
| 56-55-3  | Benzo(a)anthracene         | 330  | U |
| 218-01-9 |                            | 330  | U |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 330  | U |
| 205-99-2 | Benzo(b)fluoranthene       | 330  | U |
|          | Benzo(k)fluoranthene       | 330  | U |
| 50-32-8  | Benzo(a)pyrene             | 330  | U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene     | 330  | U |
|          | Dibenzo(a,h)anthracene     | 330  | U |
| 191-24-2 | Benzo(g,h,i)perylene       | 330  | U |

# 1D - FORM I SV-1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| CLIENT | SAMPLE | NO. |
|--------|--------|-----|
| LCS-72 | 594    |     |
|        |        |     |

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: LCS-72594      |
| Sample wt/vol:15.0 (g/mL) G               | Lab File ID: S4F4228.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: Decanted: (Y/N)               | Date Received:                |
| Concentrated Extract Volume:1000 (uL)     | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

|           |                              | CONCENTRATION UNITS:  |   |
|-----------|------------------------------|-----------------------|---|
| CAS NO.   | COMPOUND                     | (ug/L or ug/Kg) UG/KG | Q |
| 111-44-4  | Bis(2-chloroethyl)ether      | 3100                  |   |
|           | 1,3-Dichlorobenzene          | 2700                  |   |
| 106-46-7  | 1,4-Dichlorobenzene          | 2800                  |   |
| 95-50-1   | 1,2-Dichlorobenzene          | 2800                  |   |
|           | 2,2'-oxybis(1-Chloropropane) | 3000                  |   |
|           | Hexachloroethane             | 2800                  |   |
| 98-95-3   | Nitrobenzene                 | 3000                  |   |
| 78-59-1   | Isophorone                   | 2900                  |   |
|           | 1,2,4-Trichlorobenzene       | 2900                  |   |
|           | Naphthalene                  | 3000                  |   |
| 106-47-8  | 4-Chloroaniline              | 1900                  |   |
| 111-91-1  | Bis(2-chloroethoxy)methane   | 3000                  |   |
|           | Hexachlorobutadiene          | 3100                  |   |
| 91-57-6   | 2-Methylnaphthalene          | 2800                  |   |
|           | Hexachlorocyclopentadiene    | 2600                  |   |
| 91-58-7   | 2-Chloronaphthalene          | 2900                  |   |
|           | 2-Nitroaniline               | 2800                  |   |
| 131-11-3  | Dimethylphthalate            | 2800                  |   |
| 208-96-8  | Acenaphthylene               | 2800                  |   |
| 606-20-2  | 2,6-Dinitrotoluene           | 2900                  |   |
|           | 3-Nitroaniline               | 2000                  |   |
|           | Acenaphthene                 | 2800                  |   |
|           | Dibenzofuran                 | 2900                  |   |
| 121-14-2  | 2,4-Dinitrotoluene           | 2700                  |   |
| 84-66-2   | Diethylphthalate             | 2900                  |   |
| 7005-72-3 | 4-Chlorophenyl-phenylether   | 3000                  |   |
| 86-73-7   | Fluorene                     | 3000                  |   |
| 100-01-6  | 4-Nitroaniline               | 2100                  |   |
| 101-55-3  | 4-Bromophenyl-phenylether    | 3000                  |   |
| 118-74-1  | Hexachlorobenzene            | 2800                  |   |
| 85-01-8   | Phenanthrene                 | 2900                  |   |
| 120-12-7  | Anthracene                   | 2800                  |   |
| 86-74-8   | Carbazole                    | 2800                  |   |
| 206-44-0  | Fluoranthene                 | 2900                  |   |
| 129-00-0  | Pyrene                       | 3000                  |   |
| 85-68-7   | Butylbenzylphthalate         | 2900                  |   |

# 1E - FORM I SV-2 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| CLIENT | SAMPLE | NO. |
|--------|--------|-----|
| LCS-72 | 594    |     |
|        |        |     |

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: LCS-72594      |
| Sample wt/vol:15.0 (g/mL) G               | Lab File ID: S4F4228.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: Decanted: (Y/N)               | Date Received:                |
| Concentrated Extract Volume:1000 (uL)     | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

| CAS NO.  | COMPOUND                   | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG | Q |
|----------|----------------------------|--|---|
| 91-94-1  | 3,3´-Dichlorobenzidine     | 2000                                       |   |
| 56-55-3  | Benzo(a)anthracene         | 2800                                       |   |
| 218-01-9 | Chrysene                   | 3000                                       |   |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 3000                                       |   |
| 205-99-2 | Benzo(b)fluoranthene       | 2800                                       |   |
| 207-08-9 | Benzo(k)fluoranthene       | 2900                                       |   |
|          | Benzo(a)pyrene             | 2700                                       |   |
| 193-39-5 | Indeno(1,2,3-cd)pyrene     | 2600                                       |   |
| 53-70-3  | Dibenzo(a,h)anthracene     | 2600                                       |   |
| 191-24-2 | Benzo(g,h,i)perylene       | 2500                                       |   |

# 1D - FORM I SV-1 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| CLIENT | SAMPLE | NO. |
|--------|--------|-----|
| LCSD-7 | 2594   |     |
|        |        |     |

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: LCSD-72594     |
| Sample wt/vol:15.0 (g/mL) G               | Lab File ID: S4F4229.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: Decanted: (Y/N)               | Date Received:                |
| Concentrated Extract Volume:1000 (uL)     | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

|           |                              | CONCENTRATION UNITS:  |   |
|-----------|------------------------------|-----------------------|---|
| CAS NO.   | COMPOUND                     | (ug/L or ug/Kg) UG/KG | Q |
| 111-44-4  | Bis(2-chloroethyl)ether      | 3000                  |   |
|           | 1,3-Dichlorobenzene          | 2600                  |   |
|           | 1,4-Dichlorobenzene          | 2700                  |   |
| 95-50-1   | 1,2-Dichlorobenzene          | 2700                  |   |
|           | 2,2'-oxybis(1-Chloropropane) | 2900                  |   |
|           | Hexachloroethane             | 2800                  |   |
| 98-95-3   | Nitrobenzene                 | 3000                  |   |
|           | Isophorone                   | 3000                  |   |
|           | 1,2,4-Trichlorobenzene       | 2700                  |   |
|           | Naphthalene                  | 3000                  |   |
|           | 4-Chloroaniline              | 2100                  |   |
|           | Bis(2-chloroethoxy)methane   | 3000                  |   |
|           | Hexachlorobutadiene          | 2900                  |   |
|           | 2-Methylnaphthalene          | 2700                  |   |
|           | Hexachlorocyclopentadiene    | 2900                  |   |
|           | 2-Chloronaphthalene          | 2800                  |   |
|           | 2-Nitroaniline               | 2800                  |   |
| 131-11-3  | Dimethylphthalate            | 2800                  |   |
|           | Acenaphthylene               | 2800                  |   |
| 606-20-2  | 2,6-Dinitrotoluene           | 3000                  |   |
|           | 3-Nitroaniline               | 2400                  |   |
|           | Acenaphthene                 | 2800                  |   |
|           | Dibenzofuran                 | 2900                  |   |
| 121-14-2  | 2,4-Dinitrotoluene           | 2900                  |   |
| 84-66-2   | Diethylphthalate             | 2900                  |   |
| 7005-72-3 | 4-Chlorophenyl-phenylether   | 2700                  |   |
| 86-73-7   | Fluorene                     | 3000                  |   |
| 100-01-6  | 4-Nitroaniline               | 2500                  |   |
| 101-55-3  | 4-Bromophenyl-phenylether    | 3000                  |   |
| 118-74-1  | Hexachlorobenzene            | 3100                  |   |
| 85-01-8   | Phenanthrene                 | 3000                  |   |
| 120-12-7  | Anthracene                   | 2900                  |   |
|           | Carbazole                    | 2900                  |   |
| 206-44-0  | Fluoranthene                 | 3000                  |   |
| 129-00-0  | Pyrene                       | 3100                  |   |
|           | Butylbenzylphthalate         | 3100                  |   |

# 1E - FORM I SV-2 SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET

| CLIENT     | SAMPLE | NO. |  |  |  |  |
|------------|--------|-----|--|--|--|--|
| LCSD-72594 |        |     |  |  |  |  |
|            |        |     |  |  |  |  |

| Lab Name: SPECTRUM ANALYTICAL, INC.       | Contract:                     |
|---|-------------------------------|
| Lab Code: MITKEM Case No.: M1089          | Mod. Ref No.: SDG No.: SM1089 |
| Matrix: (SOIL/SED/WATER) SOIL             | Lab Sample ID: LCSD-72594     |
| Sample wt/vol:15.0 (g/mL) G               | Lab File ID: S4F4229.D        |
| Level: (LOW/MED) LOW                      | Extraction: (Type) SONC       |
| % Moisture: Decanted: (Y/N)               | Date Received:                |
| Concentrated Extract Volume: 1000 (uL)    | Date Extracted: 07/05/2013    |
| Injection Volume:1.0 (uL) GPC Factor:1.00 | Date Analyzed: 07/05/2013     |
| GPC Cleanup:(Y/N) N pH:                   | Dilution Factor: 1.0          |

| CAS NO.  | COMPOUND                   | CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG | Q |
|----------|----------------------------|--|---|
| 91-94-1  | 3,3´-Dichlorobenzidine     | 2300                                       |   |
| 56-55-3  | Benzo(a)anthracene         | 2900                                       |   |
| 218-01-9 | Chrysene                   | 3000                                       |   |
| 117-81-7 | Bis(2-ethylhexyl)phthalate | 3000                                       |   |
| 205-99-2 | Benzo(b)fluoranthene       | 3100                                       |   |
| 207-08-9 | Benzo(k)fluoranthene       | 3100                                       |   |
| 50-32-8  | Benzo(a)pyrene             | 2900                                       |   |
| 193-39-5 | Indeno(1,2,3-cd)pyrene     | 2600                                       |   |
| 53-70-3  | Dibenzo(a,h)anthracene     | 2800                                       |   |
| 191-24-2 | Benzo(g,h,i)perylene       | 2700                                       |   |

#### 2K - FORM II SV-4

#### SOIL SEMIVOLATILE DEUTERATED MONITORING COMPOUND RECOVERY

Lab Name: SPECTRUM ANALYTICAL, INC. Contract:

Lab Code: MITKEM Case No.: M1089 Mod. Ref No.: SDG No.: SM1089

Level: (LOW/MED) LOW

|    | CLIENT      | SDMC1   | SDMC2   | SDMC3   | TOT |
|----|-------------|---------|---------|---------|-----|
|    | SAMPLE NO.  | (NBZ) # | (FBP) # | (TPH) # | OUT |
| 01 | MB-72594    | 90      | 91      | 107     | 0   |
| 02 | LCS-72594   | 90      | 85      | 96      | 0   |
| 03 | LCSD-72594  | 87      | 83      | 95      | 0   |
| 04 | WC-2-061313 | 76      | 75      | 93      | 0   |
| 05 | WC-3-061313 | 76      | 80      | 97      | 0   |

|       |                          | QC LIMITS |
|-------|--------------------------|-----------|
| SDMC1 | (NBZ) = Nitrobenzene-d5  | (35-100)  |
| SDMC2 | (FBP) = 2-Fluorobiphenyl | (45-105)  |
| SDMC3 | (TPH) = Terphenyl-d14    | (30-125)  |

#### som13.06.03.A

 $<sup>\</sup>mbox{\tt\#}$  Column to be used to flag recovery values

<sup>\*</sup> Values outside of contract required QC limits

D DMC diluted out

# 3 - FORM III SOIL LABORATORY CONTROL SAMPLE RECOVERY

CLIENT SAMPLE NO.

LCS-72594

| Lab | Name:  | SPECTRUM | ANALYTICAL, | TNC   | Contract: |
|-----|--------|----------|-------------|-------|-----------|
| цар | manic. | DEFCIKON | MINDLITCHL, | TIVC. | Concract. |

Lab Code: MITKEM Case No.: M1089 Mod. Ref No.: SDG No.: SM1089

Lab Sample ID: LCS-72594 LCS Lot No.: A092773

Date Extracted: 07/05/2013 Date Analyzed (1): 07/05/2013

| COMPOUND   SPIKE   ADDED   CONCENTRATION   LCS *REC   *   LIMITS   REC.  | Extracted: 0//05/2013      |           | Date Alia.    | 1yzed (1). <u> </u> | 7/05/2013 | , |          |
|--|----------------------------|-----------|---------------|---------------------|-----------|---|----------|
| Bis(2-chloroethy1)ether   3333.0000   0.0000   3077.8942   92   40 - 105   1.3-Dichlorobenzene   3333.0000   0.0000   2681.2720   80   40 - 105   1.3-Dichlorobenzene   3333.0000   0.0000   2681.2720   80   40 - 105   1.3-Dichlorobenzene   3333.0000   0.0000   2802.5239   84   35 - 105   1.2-Dichlorobenzene   3333.0000   0.0000   2809.6290   84   45 - 95   2.2'-oxybis(1-Chloropropan   3333.0000   0.0000   2786.8920   84   35 - 116   1800.0000   3   |                            | SPIKE     | SAMPLE        | LCS                 |           |   | QC.      |
| REC.   | COMPOUND                   | ADDED     | CONCENTRATION | CONCENTRATION       | LCS %REC  | # |          |
| 1,3-Dichlorobenzene  |                            |           | l             |                     |           |   | REC.     |
| 1,3-Dichlorobenzene  | Bis(2-chloroethyl)ether    | 3333.0000 | 0.0000        | 3077.8942           | 92        |   | 40 - 105 |
| 1,4-Dichlorobensene  |                            |           |               |                     |           |   |          |
| 1,2-Dichlorobenzene  | 1,4-Dichlorobenzene        | 3333.0000 | 0.0000        | 2802.5239           | 84        |   |          |
| Hexachloroethane   | 1,2-Dichlorobenzene        | 3333.0000 | 0.0000        | 2809.6290           | 84        |   | 45 - 95  |
| Nitrobenzene   3333.0000   0.0000   3022.0847   91   40 - 115     Isophorone   3333.0000   0.0000   2926.9407   88   45 - 110     1,2,4-Trichlorobenzene   3333.0000   0.0000   2883.0554   87   45 - 110     Naphthalene   3333.0000   0.0000   3011.0077   90   40 - 105     4-Chloroaniline   3333.0000   0.0000   3011.0077   90   40 - 105     4-Chloroathoxy)methane   3333.0000   0.0000   2999.6795   90   45 - 110     Hexachlorobutadiene   3333.0000   0.0000   2999.6795   90   45 - 110     Hexachlorocyclopentadiene   3333.0000   0.0000   2779.9759   83   45 - 105     Hexachlorocyclopentadiene   3333.0000   0.0000   2776.8776   77   8 - 148     2-Chloronaphthalene   3333.0000   0.0000   2576.8776   77   8 - 148     2-Chloronaphthalene   3333.0000   0.0000   2788.2686   84   45 - 105     2-Nitroaniline   3333.0000   0.0000   2788.2686   84   45 - 120     Dimethylphthalate   3333.0000   0.0000   2786.2686   84   50 - 110     Acenaphthylene   3333.0000   0.0000   2786.2686   84   50 - 110     Acenaphthene   3333.0000   0.0000   2780.2686   84   50 - 110     Acenaphthene   3333.0000   0.0000   2270.1146   83   45 - 105     2,6-Dinitrotoluene   3333.0000   0.0000   2270.1146   83   45 - 105     2,6-Dinitrotoluene   3333.0000   0.0000   2270.1146   83   45 - 105     2,6-Dinitrotoluene   3333.0000   0.0000   2270.2529   84   45 - 110     Dibenzofuran   3333.0000   0.0000   2272.0529   84   45 - 110     Dibenzofuran   3333.0000   0.0000   2742.6559   82   50 - 115     4-Chlorophenyl-phenylether   3333.0000   0.0000   2742.6559   82   50 - 115     4-Nitroaniline   3333.0000   0.0000   2762.6520   83   35 - 115     4-Bramophenyl-phenylether   3333.0000   0.0000   2759.9316   86   55 - 110     4-Nitroaniline   3333.0000   0.0000   2759.9316   86   55 - 110     4-Nitroaniline   3333.0000   0.0000   2762.6520   83   55 - 105     4-Bramophenyl-phenylether   3333.0000   0.0000   2759.938   86   55 - 115     4-Bramophenyl-phenylether   3333.0000   0.0000   2759.938   86   55 - 115     4-Bramophenyl-phenylether   3333.0000   0.0000    | 2,2'-oxybis(1-Chloropropan | 3333.0000 | 0.0000        | 2961.6692           | 89        |   | 20 - 115 |
| Isophorone   | Hexachloroethane           | 3333.0000 | 0.0000        | 2786.8920           | 84        |   | 35 - 110 |
| 1,2,4-Trichlorobenzene   | Nitrobenzene               | 3333.0000 | 0.0000        | 3022.0847           | 91        |   | 40 - 115 |
| Naphthalene  | Isophorone                 | 3333.0000 | 0.0000        | 2926.9407           | 88        |   | 45 - 110 |
| ## Chloroaniline   3333.0000   0.0000   1872.8526   56   10 - 100   ## Bis (2-chloroethoxy) methane   3333.0000   0.0000   2999.6795   90   45 - 110   ## Cachloroethoxy) methane   3333.0000   0.0000   3132.6485   94   40 - 115   ## Cachloroethoxy methane   3333.0000   0.0000   2779.9759   83   45 - 105   ## Cachlorocyclopentadiene   3333.0000   0.0000   2576.8776   77   8 - 148   ## Cachlorocyclopentadiene   3333.0000   0.0000   2576.8776   77   8 - 148   ## Cachloronaphthalene   3333.0000   0.0000   2788.9635   84   45 - 105   ## Cachloronaphthalene   3333.0000   0.0000   2788.9635   84   45 - 120   ## Dimethylphthalate   3333.0000   0.0000   2786.2680   84   50 - 110   ## Cachlorophthylene   3333.0000   0.0000   2786.2680   84   50 - 110   ## Cachlorophthylene   3333.0000   0.0000   2786.2680   86   50 - 110   ## Cachlorophthene   3333.0000   0.0000   2247.0275   61   25 - 110   ## Cachlorophthene   3333.0000   0.0000   2792.0529   84   45 - 110   ## Dibenzofuran   3333.0000   0.0000   2792.0529   84   45 - 110   ## Cachlorophenyl-phenylether   3333.0000   0.0000   2742.6559   82   50 - 115   ## Cachlorophenyl-phenylether   3333.0000   0.0000   2242.6559   82   50 - 115   ## Cachlorophenyl-phenylether   3333.0000   0.0000   2366.5473   86   50 - 115   ## Hexachlorobenzene   3333.0000   0.0000   2366.5473   86   50 - 115   ## Cachlorophenyl-phenylether   3333.0000   0.0000   2366.5473   86   50 - 115   ## Hexachlorobenzene   3333.0000   0.0000   2366.5402   89   45 - 125   ## Hexachlorobenzene   3333.0000   0.0000   2365.3402   89   45 - 125   ## Hexachlorobenzene   3333.0000   0.0000   2365.3402   89   45 - 125   ## Butylbenzylphthalate   3333.0000   0.0000   2365.3402   89   45 - 125   ## Butylbenzylphthalate   3333.0000   0.0000   2365.3402   89   45 - 125   ## Hexachlorobenzene   3333.0000   0.0000   2365.3402   89   45 - 125   ## Butylbenzylphthalate   3333.0000   0.0000   2365.3402   89   45 - 125   ## Butylbenzylphthalate   3333.0000   0.0000   2365.3402   87   50 - 110   ## Hexachlorobenzidine   3333.0 | 1,2,4-Trichlorobenzene     | 3333.0000 | 0.0000        | 2883.0564           | 87        |   | 45 - 110 |
| Bis(2-chloroethoxy)methane   | Naphthalene                | 3333.0000 | 0.0000        | 3011.0077           | 90        |   | 40 - 105 |
| Hexachlorobutadiene  | 4-Chloroaniline            | 3333.0000 | 0.0000        | 1872.8526           | 56        |   | 10 - 100 |
| 2-Methylnaphthalene  | Bis(2-chloroethoxy)methane | 3333.0000 | 0.0000        | 2999.6795           | 90        |   | 45 - 110 |
| Hexachlorocyclopentadiene  | Hexachlorobutadiene        | 3333.0000 | 0.0000        | 3132.6485           | 94        |   | 40 - 115 |
| 2-Chloronaphthalene 3333.0000 0.0000 2851.7218 86 45 - 105 2-Mitroaniline 3333.0000 0.0000 2798.9635 84 45 - 120 Dimethylphthalate 3333.0000 0.0000 2786.6800 84 50 - 110 Acenaphthylene 3333.0000 0.0000 2750.1146 83 45 - 105 2.6-Dinitrotoluene 3333.0000 0.0000 2882.9360 86 50 - 110 3-Nitroaniline 3333.0000 0.0000 2047.0275 61 25 - 110 Acenaphthene 3333.0000 0.0000 2792.0529 84 45 - 110 Dibenzofuran 3333.0000 0.0000 2792.0529 84 45 - 110 Dibenzofuran 3333.0000 0.0000 2742.6559 82 50 - 115 Diethylphthalate 3333.0000 0.0000 2742.6559 82 50 - 115 Diethylphthalate 3333.0000 0.0000 2742.6559 82 50 - 115 Fluorene 3333.0000 0.0000 2950.1166 89 45 - 110 Fluorene 3333.0000 0.0000 2950.1166 89 45 - 110 Fluorene 3333.0000 0.0000 2950.1166 89 45 - 110 Fluorene 3333.0000 0.0000 2960.2402 89 45 - 115 Hexachlorobenzene 3333.0000 0.0000 2965.2402 89 45 - 115 Hexachlorobenzene 3333.0000 0.0000 2767.9201 83 45 - 120 Phenanthrene 3333.0000 0.0000 2767.9201 83 45 - 120 Phenanthrene 3333.0000 0.0000 2766.2402 89 45 - 115 Fluoranthene 3333.0000 0.0000 2766.9315 84 45 - 125 Butylbenzylphthalate 3333.0000 0.0000 2884.0520 87 50 - 110 Anthracene 3333.0000 0.0000 2884.0520 87 50 - 125 Butylbenzylphthalate 3333.0000 0.0000 2884.0520 87 50 - 125 Butylbenzylphthalate 3333.0000 0.0000 2782.4392 83 50 - 110 Chrysene 3333.0000 0.0000 2782.4392 83 50 - 110 Chrysene 3333.0000 0.0000 2782.4392 83 50 - 110 Bis(2-ethylhexyl)phthalate 3333.0000 0.0000 2782.4392 83 50 - 110 Bis(2-ethylhexyl)phthalate 3333.0000 0.0000 2884.0520 87 50 - 125 Benzo(b)fluoranthene 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(b)fluoranthene 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(b)fluoranthene 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(a)pyrene 3333.0000 0.0000 2876.6565 86 45 - 125  | 2-Methylnaphthalene        | 3333.0000 | 0.0000        | 2779.9759           | 83        |   | 45 - 105 |
| 2-Nitroaniline   | Hexachlorocyclopentadiene  | 3333.0000 | 0.0000        | 2576.8776           | 77        |   | 8 - 148  |
| Dimethylphthalate  | 2-Chloronaphthalene        | 3333.0000 | 0.0000        | 2851.7218           | 86        |   | 45 - 105 |
| Acenaphthylene 3333.0000 0.0000 2750.1146 83 45 - 105 2,6-Dinitrotoluene 3333.0000 0.0000 2882.9360 86 50 - 110 3-Nitroaniline 3333.0000 0.0000 2047.0275 61 25 - 110 Acenaphthene 3333.0000 0.0000 2792.0529 84 45 - 110 Dibenzofuran 3333.0000 0.0000 2792.0529 84 45 - 110 Dibenzofuran 3333.0000 0.0000 2792.0529 84 45 - 110 Dibenzofuran 3333.0000 0.0000 2742.6559 82 50 - 115 Diethylphthalate 3333.0000 0.0000 2866.5473 86 50 - 115 Diethylphthalate 3333.0000 0.0000 2866.5473 86 50 - 115 Diethylphthalate 3333.0000 0.0000 2866.5473 86 50 - 115 Diethylphthalate 3333.0000 0.0000 28950.1166 89 45 - 110 Fluorene 3333.0000 0.0000 2107.9598 63 35 - 115 Diethylphthalate 3333.0000 0.0000 2107.9598 63 35 - 115 Diethylphthalate 3333.0000 0.0000 2965.2402 89 45 - 115 Diethylphthenyl-phenylether 3333.0000 0.0000 2757.9201 83 45 - 120 Phenanthrene 3333.0000 0.0000 2757.9201 83 45 - 120 Phenanthrene 3333.0000 0.0000 2762.6520 83 55 - 105 Diethylphthalate 3333.0000 0.0000 2762.6520 83 55 - 105 Diethylphthalate 3333.0000 0.0000 2762.6520 83 55 - 105 Diethylphthalate 3333.0000 0.0000 2884.0520 87 50 - 110 Diethylphthalate 3333.0000 0.0000 2884.0520 87 50 - 115 Diethylphthalate 3333.0000 0.0000 2884.0520 87 50 - 125 Diethylphthalate 3333.0000 0.0000 2782.4392 83 50 - 110 Diethylphylphthalate 3333.0000 0.0000 2782.4392 83 5 | 2-Nitroaniline             | 3333.0000 | 0.0000        | 2798.9635           | 84        |   | 45 - 120 |
| 2,6-Dinitrotoluene   | Dimethylphthalate          | 3333.0000 | 0.0000        | 2786.2680           | 84        |   | 50 - 110 |
| 3-Nitroaniline   3333.0000   0.0000   2047.0275   61   25 - 110  | Acenaphthylene             | 3333.0000 | 0.0000        | 2750.1146           | 83        |   | 45 - 105 |
| Acenaphthene 3333.0000 0.0000 2792.0529 84 45 - 110 Dibenzofuran 3333.0000 0.0000 2933.9523 88 50 - 105 2,4-Dinitrotoluene 3333.0000 0.0000 2742.6559 82 50 - 115 Diethylphthalate 3333.0000 0.0000 2866.5473 86 50 - 115 4-Chlorophenyl-phenylether 3333.0000 0.0000 2950.1166 89 45 - 110 Fluorene 3333.0000 0.0000 308.3332 90 50 - 110 4-Nitroaniline 3333.0000 0.0000 2950.1166 89 45 - 110 4-Nitroaniline 3333.0000 0.0000 2107.9598 63 35 - 115 4-Bromophenyl-phenylether 3333.0000 0.0000 2217.9598 63 35 - 115 Hexachlorobenzene 3333.0000 0.0000 2757.9201 83 45 - 120 Phenanthrene 3333.0000 0.0000 2757.9201 83 45 - 120 Phenanthrene 3333.0000 0.0000 2757.9201 83 45 - 120 Phenanthrene 3333.0000 0.0000 2762.6520 83 55 - 105 Carbazole 3333.0000 0.0000 2762.6520 83 55 - 105 Fluoranthene 3333.0000 0.0000 2769.9315 84 45 - 115 Fluoranthene 3333.0000 0.0000 2796.9315 84 45 - 115 Pyrene 3333.0000 0.0000 2871.5938 86 55 - 115 Pyrene 3333.0000 0.0000 2871.5938 86 55 - 115 Butylbenzylphthalate 3333.0000 0.0000 2884.0520 87 50 - 125 Butylbenzylphthalate 3333.0000 0.0000 2884.0520 87 50 - 125 Butylbenzylphthalate 3333.0000 0.0000 2782.4392 83 50 - 110 Chrysene 3333.0000 0.0000 2782.4392 83 50 - 110 Bis(2-ethylhexyl)phthalate 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(b)fluoranthene 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(b)fluoranthene 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(k)fluoranthene 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(a)pyrene   | 2,6-Dinitrotoluene         | 3333.0000 | 0.0000        | 2882.9360           | 86        |   | 50 - 110 |
| Dibenzofuran   3333.0000   0.0000   2933.9523   88   50 - 105   2,4-Dinitrotoluene   3333.0000   0.0000   2742.6559   82   50 - 115   Diethylphthalate   3333.0000   0.0000   2866.5473   86   50 - 115   4-Chlorophenyl-phenylether   3333.0000   0.0000   2950.1166   89   45 - 110   Fluorene   3333.0000   0.0000   3008.3332   90   50 - 110   4-Nitroaniline   3333.0000   0.0000   2107.9598   63   35 - 115   4-Bromophenyl-phenylether   3333.0000   0.0000   2965.2402   89   45 - 115   Hexachlorobenzene   3333.0000   0.0000   2757.9201   83   45 - 125   4-Dinitroaniline   3333.0000   0.0000   2762.6520   83   55 - 105   4-Dinitroaniline   3333.0000   0.0000   2762.6520   83   55 - 105   4-Dinitroaniline   3333.0000   0.0000   2762.6520   83   55 - 105   4-Dinitroaniline   3333.0000   0.0000   2762.6520   83   55 - 115   4-Dinitroaniline   3333.0000   0.0000   2866.1324   89   45 - 125   45   45   45   45   45   45   45   | 3-Nitroaniline             | 3333.0000 | 0.0000        | 2047.0275           | 61        |   | 25 - 110 |
| 2,4-Dinitrotoluene       3333.0000       0.0000       2742.6559       82       50 - 115         Diethylphthalate       3333.0000       0.0000       2866.5473       86       50 - 115         4-Chlorophenyl-phenylether       3333.0000       0.0000       2950.1166       89       45 - 110         Fluorene       3333.0000       0.0000       3008.3332       90       50 - 110         4-Nitroaniline       3333.0000       0.0000       2107.9598       63       35 - 115         4-Bromophenyl-phenylether       3333.0000       0.0000       2965.2402       89       45 - 115         Hexachlorobenzene       3333.0000       0.0000       2757.9201       83       45 - 120         Phenanthrene       3333.0000       0.0000       2757.9201       83       45 - 120         Anthracene       3333.0000       0.0000       2762.6520       83       55 - 105         Carbazole       3333.0000       0.0000       2762.6520       83       55 - 115         Fluoranthene       3333.0000       0.0000       2871.5938       86       55 - 115         Pyrene       3333.0000       0.0000       2861.324       89       45 - 125         Butylbenzylphthalate       3333.0000   | Acenaphthene               | 3333.0000 | 0.0000        | 2792.0529           | 84        |   | 45 - 110 |
| Diethylphthalate         3333.0000         0.0000         2866.5473         86         50 - 115           4-Chlorophenyl-phenylether         3333.0000         0.0000         2950.1166         89         45 - 110           Fluorene         3333.0000         0.0000         3008.3332         90         50 - 110           4-Nitroaniline         3333.0000         0.0000         2107.9598         63         35 - 115           4-Bromophenyl-phenylether         3333.0000         0.0000         2965.2402         89         45 - 115           Hexachlorobenzene         3333.0000         0.0000         2757.9201         83         45 - 120           Phenanthrene         3333.0000         0.0000         2895.7959         87         50 - 110           Anthracene         3333.0000         0.0000         2762.6520         83         55 - 105           Carbazole         3333.0000         0.0000         2769.9315         84         45 - 115           Fluoranthene         3333.0000         0.0000         2871.5938         86         55 - 115           Pyrene         3333.0000         0.0000         2861.324         89         45 - 125           Butylbenzylphthalate         3333.0000         0.0000         2884.0520  | Dibenzofuran               | 3333.0000 | 0.0000        | 2933.9523           | 88        |   | 50 - 105 |
| 4-Chlorophenyl-phenylether       3333.0000       0.0000       2950.1166       89       45 - 110         Fluorene       3333.0000       0.0000       3008.3332       90       50 - 110         4-Nitroaniline       3333.0000       0.0000       2107.9598       63       35 - 115         4-Bromophenyl-phenylether       3333.0000       0.0000       2965.2402       89       45 - 115         Hexachlorobenzene       3333.0000       0.0000       2757.9201       83       45 - 120         Phenanthrene       3333.0000       0.0000       2895.7959       87       50 - 110         Anthracene       3333.0000       0.0000       2762.6520       83       55 - 105         Carbazole       3333.0000       0.0000       2796.9315       84       45 - 115         Fluoranthene       3333.0000       0.0000       2871.5938       86       55 - 115         Pyrene       3333.0000       0.0000       2841.5938       86       55 - 115         Butylbenzylphthalate       3333.0000       0.0000       2884.0520       87       50 - 125         Benzo(a)anthracene       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0  | 2,4-Dinitrotoluene         | 3333.0000 | 0.0000        | 2742.6559           | 82        |   | 50 - 115 |
| Fluorene         3333.0000         0.0000         3008.3332         90         50 - 110           4-Nitroaniline         3333.0000         0.0000         2107.9598         63         35 - 115           4-Bromophenyl-phenylether         3333.0000         0.0000         2965.2402         89         45 - 115           Hexachlorobenzene         3333.0000         0.0000         2757.9201         83         45 - 120           Phenanthrene         3333.0000         0.0000         2895.7959         87         50 - 110           Anthracene         3333.0000         0.0000         2762.6520         83         55 - 105           Carbazole         3333.0000         0.0000         2796.9315         84         45 - 115           Fluoranthene         3333.0000         0.0000         2871.5938         86         55 - 115           Pyrene         3333.0000         0.0000         2866.1324         89         45 - 125           Butylbenzylphthalate         3333.0000         0.0000         2884.0520         87         50 - 125           3,3´-Dichlorobenzidine         3333.0000         0.0000         2823.9874         61         10 - 130           Benzo(a)anthracene         3333.0000         0.0000         2782.4392<  | Diethylphthalate           | 3333.0000 | 0.0000        | 2866.5473           | 86        |   | 50 - 115 |
| 4-Nitroaniline       3333.0000       0.0000       2107.9598       63       35 - 115         4-Bromophenyl-phenylether       3333.0000       0.0000       2965.2402       89       45 - 115         Hexachlorobenzene       3333.0000       0.0000       2757.9201       83       45 - 120         Phenanthrene       3333.0000       0.0000       2895.7959       87       50 - 110         Anthracene       3333.0000       0.0000       2762.6520       83       55 - 105         Carbazole       3333.0000       0.0000       2796.9315       84       45 - 115         Fluoranthene       3333.0000       0.0000       2871.5938       86       55 - 115         Pyrene       3333.0000       0.0000       2966.1324       89       45 - 125         Butylbenzylphthalate       3333.0000       0.0000       2884.0520       87       50 - 125         3,3´-Dichlorobenzidine       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0.0000       2782.4392       83       50 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000<   | 4-Chlorophenyl-phenylether | 3333.0000 | 0.0000        | 2950.1166           | 89        |   | 45 - 110 |
| 4-Bromophenyl-phenylether3333.00000.00002965.24028945 - 115Hexachlorobenzene3333.00000.00002757.92018345 - 120Phenanthrene3333.00000.00002895.79598750 - 110Anthracene3333.00000.00002762.65208355 - 105Carbazole3333.00000.00002796.93158445 - 115Fluoranthene3333.00000.00002871.59388655 - 115Pyrene3333.00000.00002966.13248945 - 125Butylbenzylphthalate3333.00000.00002884.05208750 - 1253,3'-Dichlorobenzidine3333.00000.00002023.98746110 - 130Benzo(a)anthracene3333.00000.00002782.43928350 - 110Chrysene3333.00000.00003016.96359155 - 110Bis(2-ethylhexyl)phthalate3333.00000.00002950.04988945 - 125Benzo(b)fluoranthene3333.00000.00002876.65658645 - 125Benzo(k)fluoranthene3333.00000.00002876.65658645 - 125Benzo(a)pyrene3333.00000.00002676.94228050 - 110  | Fluorene                   | 3333.0000 | 0.0000        | 3008.3332           | 90        |   |          |
| Hexachlorobenzene         3333.0000         0.0000         2757.9201         83         45 - 120           Phenanthrene         3333.0000         0.0000         2895.7959         87         50 - 110           Anthracene         3333.0000         0.0000         2762.6520         83         55 - 105           Carbazole         3333.0000         0.0000         2796.9315         84         45 - 115           Fluoranthene         3333.0000         0.0000         2871.5938         86         55 - 115           Pyrene         3333.0000         0.0000         2966.1324         89         45 - 125           Butylbenzylphthalate         3333.0000         0.0000         2884.0520         87         50 - 125           3,3'-Dichlorobenzidine         3333.0000         0.0000         2023.9874         61         10 - 130           Benzo(a)anthracene         3333.0000         0.0000         2782.4392         83         50 - 110           Chrysene         3333.0000         0.0000         2950.0498         89         45 - 125           Benzo(b)fluoranthene         3333.0000         0.0000         2876.6565         86         45 - 125           Benzo(a)pyrene         3333.0000         0.0000         2676.9422  | 4-Nitroaniline             | 3333.0000 | 0.0000        | 2107.9598           | 63        |   | 35 - 115 |
| Phenanthrene         3333.0000         0.0000         2895.7959         87         50 - 110           Anthracene         3333.0000         0.0000         2762.6520         83         55 - 105           Carbazole         3333.0000         0.0000         2796.9315         84         45 - 115           Fluoranthene         3333.0000         0.0000         2871.5938         86         55 - 115           Pyrene         3333.0000         0.0000         2966.1324         89         45 - 125           Butylbenzylphthalate         3333.0000         0.0000         2884.0520         87         50 - 125           3,3'-Dichlorobenzidine         3333.0000         0.0000         2023.9874         61         10 - 130           Benzo(a)anthracene         3333.0000         0.0000         2782.4392         83         50 - 110           Chrysene         3333.0000         0.0000         3016.9635         91         55 - 110           Bis(2-ethylhexyl)phthalate         3333.0000         0.0000         2895.0498         89         45 - 125           Benzo(b)fluoranthene         3333.0000         0.0000         2876.6565         86         45 - 125           Benzo(a)pyrene         3333.0000         0.0000         2676.9  | 4-Bromophenyl-phenylether  | 3333.0000 | 0.0000        | 2965.2402           | 89        |   |          |
| Anthracene 3333.0000 0.0000 2762.6520 83 55 - 105 Carbazole 3333.0000 0.0000 2796.9315 84 45 - 115 Fluoranthene 3333.0000 0.0000 2871.5938 86 55 - 115 Pyrene 3333.0000 0.0000 2966.1324 89 45 - 125 Butylbenzylphthalate 3333.0000 0.0000 2884.0520 87 50 - 125 3,3´-Dichlorobenzidine 3333.0000 0.0000 2023.9874 61 10 - 130 Benzo(a)anthracene 3333.0000 0.0000 2782.4392 83 50 - 110 Chrysene 3333.0000 0.0000 3016.9635 91 55 - 110 Bis(2-ethylhexyl)phthalate 3333.0000 0.0000 2950.0498 89 45 - 125 Benzo(b)fluoranthene 3333.0000 0.0000 2809.4036 84 45 - 115 Benzo(k)fluoranthene 3333.0000 0.0000 2876.6565 86 45 - 125 Benzo(a)pyrene 3333.0000 0.0000 2676.9422 80 50 - 110   | Hexachlorobenzene          | 3333.0000 | 0.0000        | 2757.9201           | 83        |   |          |
| Carbazole       3333.0000       0.0000       2796.9315       84       45 - 115         Fluoranthene       3333.0000       0.0000       2871.5938       86       55 - 115         Pyrene       3333.0000       0.0000       2966.1324       89       45 - 125         Butylbenzylphthalate       3333.0000       0.0000       2884.0520       87       50 - 125         3,3´-Dichlorobenzidine       3333.0000       0.0000       2023.9874       61       10 - 130         Benzo(a)anthracene       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0.0000       3016.9635       91       55 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110   | Phenanthrene               | 3333.0000 | 0.0000        | 2895.7959           | 87        |   |          |
| Fluoranthene       3333.0000       0.0000       2871.5938       86       55 - 115         Pyrene       3333.0000       0.0000       2966.1324       89       45 - 125         Butylbenzylphthalate       3333.0000       0.0000       2884.0520       87       50 - 125         3,3'-Dichlorobenzidine       3333.0000       0.0000       2023.9874       61       10 - 130         Benzo(a)anthracene       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0.0000       3016.9635       91       55 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110  | Anthracene                 | 3333.0000 | 0.0000        | 2762.6520           | 83        |   |          |
| Pyrene       3333.0000       0.0000       2966.1324       89       45 - 125         Butylbenzylphthalate       3333.0000       0.0000       2884.0520       87       50 - 125         3,3'-Dichlorobenzidine       3333.0000       0.0000       2023.9874       61       10 - 130         Benzo(a)anthracene       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0.0000       3016.9635       91       55 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110  | Carbazole                  |           |               |                     |           |   | 45 - 115 |
| Butylbenzylphthalate       3333.0000       0.0000       2884.0520       87       50 - 125         3,3'-Dichlorobenzidine       3333.0000       0.0000       2023.9874       61       10 - 130         Benzo(a)anthracene       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0.0000       3016.9635       91       55 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110  | Fluoranthene               | 3333.0000 | 0.0000        | 2871.5938           | 86        |   | 55 - 115 |
| 3,3'-Dichlorobenzidine       3333.0000       0.0000       2023.9874       61       10 - 130         Benzo(a)anthracene       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0.0000       3016.9635       91       55 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110  | Pyrene                     | 3333.0000 | 0.0000        | 2966.1324           | 89        |   |          |
| Benzo(a) anthracene       3333.0000       0.0000       2782.4392       83       50 - 110         Chrysene       3333.0000       0.0000       3016.9635       91       55 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110   |                            | 3333.0000 |               |                     | 87        |   | 50 - 125 |
| Chrysene       3333.0000       0.0000       3016.9635       91       55 - 110         Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110  |                            |           |               |                     | 61        |   |          |
| Bis(2-ethylhexyl)phthalate       3333.0000       0.0000       2950.0498       89       45 - 125         Benzo(b)fluoranthene       3333.0000       0.0000       2809.4036       84       45 - 115         Benzo(k)fluoranthene       3333.0000       0.0000       2876.6565       86       45 - 125         Benzo(a)pyrene       3333.0000       0.0000       2676.9422       80       50 - 110  | ` '                        |           |               |                     |           |   |          |
| Benzo(b)fluoranthene         3333.0000         0.0000         2809.4036         84         45 - 115           Benzo(k)fluoranthene         3333.0000         0.0000         2876.6565         86         45 - 125           Benzo(a)pyrene         3333.0000         0.0000         2676.9422         80         50 - 110  |                            |           |               |                     |           |   |          |
| Benzo(k)fluoranthene         3333.0000         0.0000         2876.6565         86         45 - 125           Benzo(a)pyrene         3333.0000         0.0000         2676.9422         80         50 - 110  |                            |           |               |                     |           |   |          |
| Benzo(a)pyrene 3333.0000 0.0000 2676.9422 80 50 - 110  |                            |           |               |                     |           |   |          |
|  |                            |           |               |                     |           |   |          |
| Indeno(1,2,3-cd)pyrene 3333.0000 0.0000 2551.9677 77 40 - 120  |                            |           |               |                     |           |   |          |
|  | Indeno(1,2,3-cd)pyrene     | 3333.0000 | 0.0000        | 2551.9677           | 77        |   | 40 - 120 |

# 3 - FORM III SOIL LABORATORY CONTROL SAMPLE RECOVERY

CLIENT SAMPLE NO.
LCS-72594

| Lab N | Tame:  | SPECT  | RUM ANA  | LYTICAL | , INC.         | Contract       | :             |            |     |          |
|-------|--------|--------|----------|---------|----------------|----------------|---------------|------------|-----|----------|
| Lab C | ode:   | MITKE  | М        | Case N  | Jo.: M1089     | Mod. Ref       | No.:          | SDG        | No. | : SM1089 |
| Lab S | ample  | ID:    | LCS-72   | 594     |                | LCS Lot        | No.: A092     | 773        |     |          |
| Date  | Extra  | cted:  | 07/05/   | 2013    |                | Date Ana       | lyzed (1): (  | 07/05/2013 | }   |          |
|       |        |        |          |         | SPIKE          | SAMPLE         | LCS           |            |     | QC.      |
|       |        | COM    | IPOUND   |         | ADDED          | CONCENTRATION  | CONCENTRATION | LCS %REC   | #   | LIMITS   |
|       |        |        |          |         |                |                |               |            |     | REC.     |
|       | Diben  | zo(a,h | )anthrac | ene     | 3333.0000      | 0.0000         | 2553.3208     | 3 77       |     | 40 - 125 |
|       | Benzo  | (g,h,i | )perylen | ıe      | 3333.0000      | 0.0000         | 2535.0933     | 76         |     | 40 - 125 |
|       |        |        | ed to fl |         | very and RPD v | values with an | asterisk      |            |     |          |
|       |        |        | ~        |         |                |                |               |            |     |          |
| Spike | Recove | ery:   | 0 oı     | ut of _ | 46 outside     | limits         |               |            |     |          |
|       |        |        |          |         |                |                |               |            |     |          |
| COMME | мтс:   |        |          |         |                |                |               |            |     |          |

SW846

# 3 - FORM III SOIL LABORATORY CONTROL SAMPLE DUPLICATE RECOVERY

EPA SAMPLE NO.

LCSD-72594

Lab Name: SPECTRUM ANALYTICAL, INC. Contract:

Lab Code: MITKEM Case No.: M1089 Mod. Ref No.: SDG No.: SM1089

Lab Sample ID: LCSD-72594 LCS Lot No.: A092773

|                            | SPIKE     | LCSD          |           |   |      |   | QC  | LIMITS   |
|----------------------------|-----------|---------------|-----------|---|------|---|-----|----------|
|                            | ADDED     | CONCENTRATION | LCSD %REC | # | %RPD | # |     |          |
| COMPOUND                   |           |               |           |   |      |   | RPD | REC.     |
| Bis(2-chloroethyl)ether    | 3333.0000 |               |           |   | 3    |   | 40  | 40 - 105 |
| 1,3-Dichlorobenzene        | 3333.0000 |               |           |   | 4    |   | 40  | 40 - 100 |
| 1,4-Dichlorobenzene        | 3333.0000 |               |           |   | 4    |   | 40  | 35 - 105 |
| 1,2-Dichlorobenzene        | 3333.0000 |               |           |   | 4    |   | 40  | 45 - 95  |
| 2,2'-oxybis(1-Chloropropan | 3333.0000 |               |           |   | 2    |   | 40  | 20 - 115 |
| Hexachloroethane           | 3333.0000 |               |           |   | 0    |   | 40  | 35 - 110 |
| Nitrobenzene               | 3333.0000 |               |           |   | 0    |   | 40  | 40 - 115 |
| Isophorone                 | 3333.0000 |               |           |   | 3    |   | 40  | 45 - 110 |
| 1,2,4-Trichlorobenzene     | 3333.0000 |               | 81        |   | 7    |   | 40  | 45 - 110 |
| Naphthalene                | 3333.0000 |               |           |   | 1    |   | 40  | 40 - 105 |
| 4-Chloroaniline            | 3333.0000 | 2061.6653     | 62        |   | 10   |   | 40  | 10 - 100 |
| Bis(2-chloroethoxy)methane | 3333.0000 | 3036.1012     | 91        |   | 1    |   | 40  | 45 - 110 |
| Hexachlorobutadiene        | 3333.0000 | 2881.9734     | 86        |   | 9    |   | 40  | 40 - 115 |
| 2-Methylnaphthalene        | 3333.0000 | 2705.3865     | 81        |   | 2    |   | 40  | 45 - 105 |
| Hexachlorocyclopentadiene  | 3333.0000 | 2875.4951     | 86        |   | 11   |   | 40  | 8 - 148  |
| 2-Chloronaphthalene        | 3333.0000 | 2795.5240     | 84        |   | 2    |   | 40  | 45 - 105 |
| 2-Nitroaniline             | 3333.0000 | 2822.8185     | 85        |   | 1    |   | 40  | 45 - 120 |
| Dimethylphthalate          | 3333.0000 | 2755.7175     | 83        |   | 1    |   | 40  | 50 - 110 |
| Acenaphthylene             | 3333.0000 | 2787.1719     | 84        |   | 1    |   | 40  | 45 - 105 |
| 2,6-Dinitrotoluene         | 3333.0000 | 2982.2556     | 89        |   | 3    |   | 40  | 50 - 110 |
| 3-Nitroaniline             | 3333.0000 | 2442.1568     | 73        |   | 18   |   | 40  | 25 - 110 |
| Acenaphthene               | 3333.0000 | 2839.8312     | 85        |   | 1    |   | 40  | 45 - 110 |
| Dibenzofuran               | 3333.0000 | 2933.7094     | 88        |   | 0    |   | 40  | 50 - 105 |
| 2,4-Dinitrotoluene         | 3333.0000 | 2948.4476     | 88        |   | 7    |   | 40  | 50 - 115 |
| Diethylphthalate           | 3333.0000 | 2934.4081     | 88        |   | 2    |   | 40  | 50 - 115 |
| 4-Chlorophenyl-phenylether | 3333.0000 | 2748.2643     | 82        |   | 8    |   | 40  | 45 - 110 |
| Fluorene                   | 3333.0000 | 2956.1481     | 89        |   | 1    |   | 40  | 50 - 110 |
| 4-Nitroaniline             | 3333.0000 | 2504.8490     | 75        |   | 17   |   | 40  | 35 - 115 |
| 4-Bromophenyl-phenylether  | 3333.0000 | 3042.2561     | 91        |   | 2    |   | 40  | 45 - 115 |
| Hexachlorobenzene          | 3333.0000 | 3093.2336     | 93        |   | 11   |   | 40  | 45 - 120 |
| Phenanthrene               | 3333.0000 | 2953.7852     | 89        |   | 2    |   | 40  | 50 - 110 |
| Anthracene                 | 3333.0000 | 2870.4793     | 86        |   | 4    |   | 40  | 55 - 105 |
| Carbazole                  | 3333.0000 | 2935.6265     | 88        |   | 5    |   | 40  | 45 - 115 |
| Fluoranthene               | 3333.0000 | 3004.9434     | 90        |   | 5    |   | 40  | 55 - 115 |
| Pyrene                     | 3333.0000 | 3084.2506     | 93        |   | 4    |   | 40  | 45 - 125 |
| Butylbenzylphthalate       | 3333.0000 | 3072.3630     | 92        |   | 6    |   | 40  | 50 - 125 |
| 3,3´-Dichlorobenzidine     | 3333.0000 | 2288.5234     | 69        |   | 12   |   | 40  | 10 - 130 |
| Benzo(a)anthracene         | 3333.0000 | 2921.8288     | 88        |   | 6    |   | 40  | 50 - 110 |
| Chrysene                   | 3333.0000 | 3047.8924     | 91        |   | 0    |   | 40  | 55 - 110 |
| Bis(2-ethylhexyl)phthalate | 3333.0000 | 3034.4535     | 91        |   | 2    |   | 40  | 45 - 125 |
| Benzo(b)fluoranthene       | 3333.0000 |               |           |   | 9    |   | 40  | 45 - 115 |
| Benzo(k)fluoranthene       | 3333.0000 | 3128.2197     | 94        |   | 9    |   | 40  | 45 - 125 |
| Benzo(a)pyrene             | 3333.0000 |               |           |   | 7    |   | 40  | 50 - 110 |
| Indeno(1,2,3-cd)pyrene     | 3333.0000 |               |           |   | 0    |   | 40  | 40 - 120 |
| Dibenzo(a,h)anthracene     | 3333.0000 |               |           |   | 8    |   | 40  | 40 - 125 |
| Benzo(g,h,i)perylene       | 3333.0000 |               |           |   | 5    |   | 40  | 40 - 125 |

# 3 - FORM III SOIL LABORATORY CONTROL SAMPLE DUPLICATE RECOVERY

EPA SAMPLE NO.
LCSD-72594

| Lab Name:                        | SPECTRUM ANALYTIC         | Contract       | :                     |           |      |        |         |        |
|----------------------------------|---------------------------|----------------|-----------------------|-----------|------|--------|---------|--------|
| Lab Code: MITKEM Case No.: M1089 |                           |                | Mod. Ref              | No.:      |      | SI     | OG No.: | SM1089 |
| Lab Sample                       | ID: <u>LCSD-72594</u>     |                | LCS Lot 1             | No.: A    | 0927 | 73     |         |        |
|                                  |                           | SPIKE<br>ADDED | LCSD<br>CONCENTRATION | LCSD %REC | # 8  | kRPD # | QC I    | IMITS  |
|                                  | COMPOUND                  |                |                       |           |      |        | RPD     | REC.   |
|                                  | be used to flag rectaints | covery and RPD | values with an        | asterisk  |      |        |         |        |
| RPD: 0                           | out of 46 outs            | ide limits     |                       |           |      |        |         |        |
| Spike Recov                      | ery: 0 out of             | 46outside      | limits                |           |      |        |         |        |
| COMMENTS:                        |                           |                |                       |           |      |        |         |        |

# 4C - FORM IV SV SEMIVOLATILE METHOD BLANK SUMMARY

CLIENT SAMPLE NO.
MB-72594

| Lab Name: SI | PECTRUM ANA  | ALYTICAL, INC.  | Contract:        |                 |
|--------------|--------------|-----------------|------------------|-----------------|
| Lab Code: MI | ITKEM        | Case No.: M1089 | Mod. Ref No.:    | SDG No.: SM1089 |
| Lab File ID: | S4F422       | 7.D             | Lab Sample ID:   | MB-72594        |
| Instrument I | D: <u>S4</u> |                 | Date Extracted:  | 07/05/2013      |
| Matrix: (SOI | L/SED/WATER  | R) SOIL         | Date Analyzed:   | 07/05/2013      |
| Level: (LOW/ | MED) LOW     |                 | Time Analyzed:   | 19:40           |
| Extraction:  | (Type) SO    | NC              | GPC Cleanup: (Y/ | N) N            |

|    | EPA         | LAB        | LAB       | DATE       |
|----|-------------|------------|-----------|------------|
|    | SAMPLE NO.  | SAMPLE ID  | FILE ID   | ANALYZED   |
| 01 | LCS-72594   | LCS-72594  | S4F4228.D | 07/05/2013 |
| 02 | LCSD-72594  | LCSD-72594 | S4F4229.D | 07/05/2013 |
| 03 | WC-2-061313 | M1089-01A  | S4F4234.D | 07/05/2013 |
| 04 | WC-3-061313 | M1089-02A  | S4F4235.D | 07/05/2013 |

| COMMENTS: |  |  |  |
|-----------|--|--|--|
|           |  |  |  |

som13.06.03.A Page 1 of 1 SW846

M1089 Page 23 of 37

#### 8C - FORM VIII SV-1

#### SEMIVOLATILE INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: SPECTRUM ANALYTICAL, INC. Contract:

Lab Code: MITKEM Case No.: M1089 Mod. Ref No.: SDG No.: SM1089

EPA Sample No.(SSTD020##) SSTD0254C Date Analyzed: 07/05/2013

Lab File ID (Standard): S4F4211.D Time Analyzed: 12:15

Instrument ID: S4

|    |             | IS1 (DCB) | IS2 (NPT) |   |        |   |       | IS3 (ANT) |        |   |       |   |
|----|-------------|-----------|-----------|---|--------|---|-------|-----------|--------|---|-------|---|
|    |             | AREA ‡    | RT        | # | AREA   | # | RT    | #         | AREA   | # | RT    | # |
|    | 12 HOUR STD | 51686     | 5.066     |   | 204769 |   | 6.267 |           | 102514 |   | 7.974 |   |
|    | UPPER LIMIT | 103372    | 5.566     |   | 409538 |   | 6.767 |           | 205028 |   | 8.474 |   |
|    | LOWER LIMIT | 25843     | 4.566     |   | 102385 |   | 5.767 |           | 51257  |   | 7.474 |   |
|    | SAMPLE NO.  |           |           |   |        |   |       |           |        |   |       |   |
| 01 | MB-72594    | 75734     | 5.066     |   | 252627 |   | 6.267 |           | 127604 |   | 7.964 |   |
| 02 | LCS-72594   | 63941     | 5.066     |   | 229536 |   | 6.267 |           | 116429 |   | 7.964 |   |
| 03 | LCSD-72594  | 73729     | 5.066     |   | 265456 |   | 6.267 |           | 134332 |   | 7.964 |   |
| 04 | WC-2-061313 | 71772     | 5.066     |   | 249244 |   | 6.266 |           | 130784 |   | 7.964 |   |
| 05 | WC-3-061313 | 72388     | 5.066     |   | 262214 |   | 6.267 |           | 135725 |   | 7.964 |   |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = 200% of internal standard area

AREA LOWER LIMIT = 50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside contract required QC limits with an asterisk.

som 13.06.03.A Page 1 of 1

SW846

#### 8D - FORM VIII SV-2

#### SEMIVOLATILE INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: SPECTRUM ANALYTICAL, INC. Contract:

Lab Code: MITKEM Case No.: M1089 Mod. Ref No.: SDG No.: SM1089

EPA Sample No.(SSTD020##) SSTD0254C Date Analyzed: 07/05/2013

Lab File ID (Standard): S4F4211.D Time Analyzed: 12:15

Instrument ID: S4 GC Column: Rxi-5sil MS ID: 0.25 (mm)

|    |             | IS4 (PHN) |   |       |   | IS5 (CRY) |   |        |   | IS6 (PRY) |   |        |   |
|----|-------------|-----------|---|-------|---|-----------|---|--------|---|-----------|---|--------|---|
|    |             | AREA      | # | RT ;  | # | AREA      | # | RT     | # | AREA      | # | RT     | # |
| -  | 12 HOUR STD | 174197    |   | 9.423 |   | 179233    |   | 12.259 |   | 158976    |   | 14.692 |   |
| ,  | UPPER LIMIT | 348394    |   | 9.923 |   | 358466    |   | 12.759 |   | 317952    |   | 15.192 |   |
| -  | LOWER LIMIT | 87099     |   | 8.923 |   | 89617     |   | 11.759 |   | 79488     |   | 14.192 |   |
|    | SAMPLE NO.  |           |   |       |   |           |   |        |   |           |   |        |   |
| 01 | MB-72594    | 217039    |   | 9.423 |   | 203772    |   | 12.104 |   | 184539    |   | 14.495 |   |
| 02 | LCS-72594   | 203063    |   | 9.424 |   | 194765    |   | 12.104 |   | 180976    |   | 14.495 |   |
| 03 | LCSD-72594  | 227001    |   | 9.423 |   | 218524    |   | 12.104 |   | 187930    |   | 14.495 |   |
| 04 | WC-2-061313 | 220846    |   | 9.423 |   | 204454    |   | 12.104 |   | 187637    |   | 14.495 |   |
| 05 | WC-3-061313 | 229976    |   | 9.423 |   | 218745    |   | 12.104 |   | 213864    |   | 14.495 |   |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = 200% of internal standard area

AREA LOWER LIMIT = 50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside contract required QC limits with an asterisk.

som13.06.03.A Page 1 of 1 SW846



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

\* Metals \*

M1089

## REPORT NARRATIVE

Spectrum Analytical, Inc. Featuring Hanibal Technology, RI Division.

Client: GZA GeoEnvironmental of NY Buffalo

**Project: Former Signore Facility** 

Laboratory Workorder / SDG #: M1089 SW846 6010C, SW846 7471B

#### I. SAMPLE RECEIPT

No exceptions or unusual conditions were encountered unless a Sample Condition Notification Form, or other record of communication is included with the Sample Receipt Documentation.

### II. HOLDING TIMES

## A. Sample Preparation:

All samples were prepared within the method-specified holding times.

### B. Sample Analysis:

All samples were analyzed within the method-specified holding times.

## III. METHODS

Samples were analyzed following procedures in laboratory test code: SW846 6010C, SW846 7471B.

#### IV. PREPARATION

Soil Samples were prepared following procedures in laboratory test code: SW3050B.

Soil Samples were prepared following procedures in laboratory test code: SW7471B.

### V. INSTRUMENTATION

The following instrumentation was used:

M1089 Page 27 of 37

Instrument Code: FIMS2 Instrument Type: CVAA

Description: FIMS

Manufacturer: Perkin-Elmer

Model: FIMS100

Instrument Code: OPTIMA3

Instrument Type: ICP

Description: Optima ICP-OES Manufacturer: Perkin-Elmer

Model: 4300 DV

## VI. ANALYSIS

## A. Calibration:

Calibrations met the method/SOP acceptance criteria.

## B. Blanks:

All method blanks were within the acceptance criteria.

## C. Spikes:

## 1. Laboratory Control Spikes (LCS):

Percent recoveries for laboratory control samples were within the QC limits.

## 2. Matrix spike (MS):

A matrix spike was not performed on any sample in this SDG.

## D. Post Digestion Spike (PDS):

A post-digestion spike was not performed on any sample in this SDG.

## E. Duplicate sample:

A duplicate analysis was not performed on any sample in this SDG.

## F. Serial Dilution (SD):

A serial dilution was not performed on any sample in this SDG.

## G. Samples:

M1089 Page 28 of 37

No other unusual occurrences were noted during sample analysis.

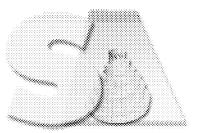
I certify that this data package is in compliance with the terms and conditions agreed to by the client and Spectrum, both technically and for completeness, except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

Shann B Law le

Signed:

Date: 07/12/13

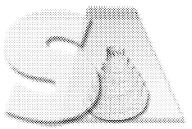
M1089 Page 29 of 37



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

# Data Flag/Qualifiers:

- U Not Detected. This compound was analyzed-for but not detected. For most analyses the reporting limit (lowest standard concentration) is the value listed. For Department of Defense programs, this is the Limit of Detection (LOD).
- J This flag indicates an estimated value due to either
  - the compound was detected below the reporting limit, or
  - estimated concentration for Tentatively Identified Compound
- B This flag indicates the compound was also detected in the associated Method Blank. The B flag has an alternative meaning for Inorganics analyses reported using CLP ILM-type metals forms, indicating a "trace" concentration below the reporting limit and equal to or above the detection limit.
- D For Organics analysis, this flag indicates the compound concentration was obtained from a secondary dilution analysis
- E This flag indicates the compound concentration exceeded the Calibration Range. The E flag has an alternative meaning for Inorganics analyses reported using CLP metals forms, indicating an estimated concentration due to the presence of interferences, as determined by the serial dilution analysis.
- P This flag is used for pesticides/PCB/herbicide compound when there is a greater than 40% difference for detected concentration between the two GC columns used for primary and confirmation analyses. This difference typically indicates an interference, causing one value to be unusually high. The **lower** of the two values is generally reported on the Form 1, and both values reported on the Form 10.
- A Used to flag semivolatile organic Tentatively Identified Compound library search results for compounds identified as aldol condensation byproducts.
- N Used to flag results for volatile and semivolatile Organics analysis Tentatively Identified Compounds where an analyte has passed the identification criteria, and is considered to be positively identified. For Inorganics analysis the N flag indicates the matrix spike recovery falls outside of the control limit.
- \* For Inorganics analysis the \* flag indicates Relative Percent Difference for duplicate analyses is outside of the control limit.



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

# Sample ID Suffixes

- DL Diluted analysis. The sample was diluted and reanalyzed. The DL may be followed by a digit if more than one diluted reanalysis is provided. The DL suffix is not attached to an analysis initially performed at dilution, only to reanalyses performed at dilution
- RE Reanalysis. Appended to the client sample ID to indicate a reextraction and reanalysis or a reanalysis of the original sample extract.
- RA Reanalysis. Appended to the laboratory sample ID indicates a reanalysis of the original sample extract.
- RX Reextraction. Appended to the laboratory sample ID indicates a reextraction of the sample.
- MS Matrix Spike.
- MSD Matrix Spike Duplicate
- DUP Duplicate analysis
- SD Serial Dilution
- PS Post-digestion or Post-distillation spike. For metals or inorganic analyses

EPA SAMPLE NO.

WC-2-061313

### INORGANIC ANALYSIS DATA SHEET

|--|

Lab Code: MITKEM Case No.: SAS No.: SDG No.: SM1089

Matrix (soil/water): SOIL Lab Sample ID: M1089-01

Level (low/med): MED Date Received: 07/03/2013

% Solids: 88.0

Concentration Units (ug/L or mg/kg dry weight): MG/KG

| CAS No.   | Analyte   | Concentration | С | Q | М  |
|-----------|-----------|---------------|---|---|----|
| 7429-90-5 | Aluminum  | 7100          |   |   | P  |
| 7440-36-0 | Antimony  | 0.37          | U |   | P  |
| 7440-38-2 | Arsenic   | 9.6           |   |   | P  |
| 7440-39-3 | Barium    | 95.5          |   |   | P  |
| 7440-41-7 | Beryllium | 0.32          |   |   | P  |
| 7440-43-9 | Cadmium   | 0.015         | U |   | P  |
| 7440-70-2 | Calcium   | 3120          |   |   | P  |
| 7440-47-3 | Chromium  | 9.9           |   |   | P  |
| 7440-48-4 | Cobalt    | 6.5           |   |   | P  |
| 7440-50-8 | Copper    | 24.1          |   |   | P  |
| 7439-89-6 | Iron      | 15400         |   |   | P  |
| 7439-92-1 | Lead      | 11.9          |   |   | P  |
| 7439-95-4 | Magnesium | 2340          |   |   | P  |
| 7439-96-5 | Manganese | 631           |   |   | Р  |
| 7439-97-6 | Mercury   | 0.012         | В |   | CV |
| 7440-02-0 | Nickel    | 14.2          |   |   | P  |
| 7440-09-7 | Potassium | 677           |   |   | P  |
| 7782-49-2 | Selenium  | 1.1           | В |   | P  |
| 7440-22-4 | Silver    | 0.063         | U |   | P  |
| 7440-23-5 | Sodium    | 53.7          |   |   | P  |
| 7440-28-0 | Thallium  | 0.22          | U |   | P  |
| 7440-62-2 | Vanadium  | 9.1           |   |   | P  |
| 7440-66-6 | Zinc      | 105           |   |   | P  |

| Commen | its: |  |  |  |
|--------|------|--|--|--|
|        |      |  |  |  |
| -      |      |  |  |  |
| -      |      |  |  |  |
|        |      |  |  |  |

M1089

EPA SAMPLE NO.

WC-3-061313

# INORGANIC ANALYSIS DATA SHEET

Lab Code: MITKEM Case No.: SAS No.: SDG No.: SM1089

Matrix (soil/water): SOIL Lab Sample ID: M1089-02

Level (low/med): MED Date Received: 07/03/2013

% Solids: 91.1

Concentration Units (ug/L or mg/kg dry weight): MG/KG

|           | _         | T             |   | _ |    |
|-----------|-----------|---------------|---|---|----|
| CAS No.   | Analyte   | Concentration | С | Q | M  |
| 7429-90-5 | Aluminum  | 8220          |   |   | P  |
| 7440-36-0 | Antimony  | 0.39          | U |   | P  |
| 7440-38-2 | Arsenic   | 10.1          |   |   | P  |
| 7440-39-3 | Barium    | 102           |   |   | P  |
| 7440-41-7 | Beryllium | 0.39          |   |   | P  |
| 7440-43-9 | Cadmium   | 0.016         | U |   | P  |
| 7440-70-2 | Calcium   | 3390          |   |   | P  |
| 7440-47-3 | Chromium  | 11.1          |   |   | P  |
| 7440-48-4 | Cobalt    | 8.1           |   |   | P  |
| 7440-50-8 | Copper    | 20.7          |   |   | P  |
| 7439-89-6 | Iron      | 18400         |   |   | P  |
| 7439-92-1 | Lead      | 12.9          |   |   | P  |
| 7439-95-4 | Magnesium | 2790          |   |   | P  |
| 7439-96-5 | Manganese | 667           |   |   | P  |
| 7439-97-6 | Mercury   | 0.013         | В |   | CV |
| 7440-02-0 | Nickel    | 16.7          |   |   | P  |
| 7440-09-7 | Potassium | 717           |   |   | P  |
| 7782-49-2 | Selenium  | 0.71          | В |   | P  |
| 7440-22-4 | Silver    | 0.11          | В |   | P  |
| 7440-23-5 | Sodium    | 51.5          | В |   | P  |
| 7440-28-0 | Thallium  | 0.23          | U |   | P  |
| 7440-62-2 | Vanadium  | 10.3          |   |   | P  |
| 7440-66-6 | Zinc      | 95.5          |   |   | P  |
| L         | 1         | l .           | L | 1 |    |

| Commen | its: |  |  |  |
|--------|------|--|--|--|
|        |      |  |  |  |
| •      |      |  |  |  |
|        |      |  |  |  |
|        |      |  |  |  |

## LABORATORY CONTROL SAMPLE

| Lab Name: Spectrum Analytical, Inc. |            | ytical, Inc. | Contract: | 21.0056491.00 |            |        |  |
|-------------------------------------|------------|--------------|-----------|---------------|------------|--------|--|
| Lab Code:                           | MITKEM     | Case No.:    | SAS No.:  |               | SDG No.:   | SM1089 |  |
| Solid LCS                           | Source:    |              |           |               | LCS(D) ID: |        |  |
| Aqueous Lo                          | CS Source: |              |           |               | LCS-72597  |        |  |

|           | Aqu  | eous (ug/I | ٦) | Solid (mg/Kg) |        |       |        |       |
|-----------|------|------------|----|---------------|--------|-------|--------|-------|
| Analyte   | True | Found      | %R | True          | Found  | C Lim | its    | %R    |
| Aluminum  |      |            |    | 455.0         | 441.8  | 364   | 546.0  | 97.1  |
| Antimony  |      |            |    | 22.8          | 23.1   | 18.2  | 27.3   | 101.3 |
| Arsenic   |      |            |    | 22.8          | 22.4   | 18.2  | 27.3   | 98.2  |
| Barium    |      |            |    | 455.0         | 463.7  | 364   | 546.0  | 101.9 |
| Beryllium |      |            |    | 11.4          | 11.2   | 9.1   | 13.6   | 98.2  |
| Cadmium   |      |            |    | 11.4          | 11.3   | 9.1   | 13.6   | 99.1  |
| Calcium   |      |            |    | 1135.0        | 1090.6 | 908   | 1362.0 | 96.1  |
| Chromium  |      |            |    | 45.5          | 45.3   | 36.4  | 54.6   | 99.6  |
| Cobalt    |      |            |    | 113.5         | 112.9  | 90.8  | 136.2  | 99.5  |
| Copper    |      |            |    | 56.5          | 56.3   | 45.2  | 67.8   | 99.6  |
| Iron      |      |            |    | 227.5         | 236.7  | 182   | 273.0  | 104.0 |
| Lead      |      |            |    | 22.8          | 22.4   | 18.2  | 27.3   | 98.2  |
| Magnesium |      |            |    | 1135.0        | 1151.0 | 908   | 1362.0 | 101.4 |
| Manganese |      |            |    | 113.5         | 113.8  | 90.8  | 136.2  | 100.3 |
| Nickel    |      |            |    | 113.5         | 112.6  | 90.8  | 136.2  | 99.2  |
| Potassium |      |            |    | 1135.0        | 1140.4 | 908   | 1362.0 | 100.5 |
| Selenium  |      |            |    | 22.8          | 21.6   | 18.2  | 27.3   | 94.7  |
| Silver    |      |            |    | 56.5          | 54.8   | 42.4  | 67.8   | 97.0  |
| Sodium    |      |            |    | 1135.0        | 1142.2 | 908   | 1362.0 | 100.6 |
| Thallium  |      |            |    | 22.8          | 21.1   | 18.2  | 27.3   | 92.5  |
| Vanadium  |      |            |    | 113.5         | 112.6  | 90.8  | 136.2  | 99.2  |
| Zinc      |      |            |    | 113.5         | 110.3  | 90.8  | 136.2  | 97.2  |

7

## LABORATORY CONTROL SAMPLE

| Lab Name:  | Spectrum Analy | tical, Inc. | Contract: | 21.0056491.00 |            |        |  |
|------------|----------------|-------------|-----------|---------------|------------|--------|--|
| Lab Code:  | MITKEM         | Case No.:   | SAS No.:  |               | SDG No.:   | SM1089 |  |
| Solid LCS  | Source:        |             |           |               | LCS(D) ID: |        |  |
| Aqueous LO | S Source:      |             |           |               | LCS-72621  |        |  |

|         | Aque | eous (ug/I | ٦) |      | Solid ( | mg/Kg) |     |      |
|---------|------|------------|----|------|---------|--------|-----|------|
| Analyte | True | Found      | %R | True | Found C | Limits |     | %R   |
| Mercury |      |            |    | 0.8  | 0.7     | 0.6    | 0.9 | 87.5 |

3

BLANKS

| Lab Name:  | Spectrum Analy  | tical, Inc. | Contract: | 21.0056491.00 |          |        |  |  |
|------------|-----------------|-------------|-----------|---------------|----------|--------|--|--|
| Lab Code:  | MITKEM          | Case No.:   | SAS No.:  | S             | SDG No.: | SM1089 |  |  |
|            |                 |             |           | Blank ID:     |          |        |  |  |
| Preparatio | on Blank Concen | /KG         | MB-7262   | 1             |          |        |  |  |

# FIMS2\_130709B

|         | Initial     |   |                        |   |   |   |  |             |       |   |    |
|---------|-------------|---|------------------------|---|---|---|--|-------------|-------|---|----|
|         | Calibration | n | Continuing Calibration |   |   |   |  | Preparation |       |   |    |
|         | Blank (ug/L | ) | Blank (ug/L)           |   |   |   |  | Blank       |       |   |    |
| Analyte |             | С | 07/09/13 14:29         | С | С | 7 |  | С           |       | С | M  |
| Mercury | 0.028       | U | 0.028                  | U |   |   |  |             | 0.002 | U | CV |

M1089 Page 36 of 37

BLANKS

| Lab Nar | me:  | Spectrum Analy | tical, Inc.        | Contract: | 21.0056491.00 |          |           |  |  |
|---------|------|----------------|--------------------|-----------|---------------|----------|-----------|--|--|
| Lab Coo | de:  | MITKEM         | Case No.:          | SAS No.:  | SDG           | No.:     | SM1089    |  |  |
| Prepara | atio | n Blank Matrix | (soil/water): SOIL | ı         |               | Method I | 3lank ID: |  |  |

Preparation Blank Concentration Units (ug/L or mg/kg): MG/KG

## OPTIMA3\_130709A

MB-72597

|           |             |     |               | <b>-</b>               | 11MA3_130/0:  |   |   |       |             |   |  |
|-----------|-------------|-----|---------------|------------------------|---------------|---|---|-------|-------------|---|--|
|           | Initial     |     |               |                        |               |   |   |       |             |   |  |
|           | Calibration | n   | Co            | Continuing Calibration |               |   |   |       | Preparation |   |  |
|           | Blank (ug/I | ر ر |               | Blank (ug/L)           |               |   |   | Blank |             |   |  |
| Analyte   |             | С   | 07/09/13 8:40 | С                      | 07/09/13 9:21 | С | C |       | С           | M |  |
| Aluminum  | 66.0        | U   | 66.0          | U                      | 66.0          | U |   | 1.200 | U           | Р |  |
| Antimony  | 9.3         | U   | 9.3           | U                      | 9.3           | U |   | 0.380 | U           | P |  |
| Arsenic   | 4.5         | В   | 4.3           | U                      | 4.3           | U |   | 0.410 | U           | P |  |
| Barium    | 1.1         | U   | 1.1           | В                      | 1.1           | U |   | 0.043 | В           | P |  |
| Beryllium | 0.3         | U   | 0.3           | U                      | 0.3           | U |   | 0.002 | U           | P |  |
| Cadmium   | 0.9         | U   | 0.9           | U                      | 0.9           | U |   | 0.015 | U           | P |  |
| Calcium   | 110.0       | U   | 110.0         | U                      | 110.0         | U |   | 6.100 | U           | P |  |
| Chromium  | 0.6         | U   | 0.6           | U                      | 0.6           | U |   | 0.019 | U           | Р |  |
| Cobalt    | 0.7         | U   | 0.7           | U                      | 0.7           | U |   | 0.044 | U           | P |  |
| Copper    | 3.6         | U   | 3.6           | U                      | 3.6           | U |   | 0.110 | U           | Р |  |
| Iron      | 31.0        | U   | 31.0          | U                      | 31.0          | U |   | 1.500 | U           | P |  |
| Lead      | 4.2         | U   | 4.2           | U                      | 4.2           | U |   | 0.170 | U           | P |  |
| Magnesium | 76.0        | U   | 76.0          | U                      | 76.0          | U |   | 0.669 | В           | P |  |
| Manganese | 10.0        | U   | 10.0          | U                      | 10.0          | U |   | 0.130 | U           | P |  |
| Nickel    | 0.9         | U   | 0.8           | U                      | 0.8           | U |   | 0.043 | U           | P |  |
| Potassium | 157.7       | В   | 76.0          | U                      | 78.0          | В |   | 3.400 | U           | P |  |
| Selenium  | 12.0        | U   | 12.0          | U                      | 12.0          | U |   | 0.640 | U           | Р |  |
| Silver    | 6.9         | U   | 6.9           | U                      | 6.9           | U |   | 0.064 | U           | Р |  |
| Sodium    | 29.0        | U   | 29.0          | U                      | 29.0          | U |   | 1.100 | U           | Р |  |
| Thallium  | 6.2         | U   | 6.2           | U                      | 6.2           | U |   | 0.220 | U           | P |  |
| Vanadium  | 1.1         | U   | 1.1           | U                      | 1.1           | U |   | 0.060 | U           | P |  |
| Zinc      | 4.9         | U   | 4.9           | U                      | 4.9           | U |   | 0.227 | В           | P |  |

M1089 Page 37 of 37



# APPENDIX C

# **Material Safety Data Sheet**

Manufacturers Name & Address

**Ecocycle Corporation** 

694-2, Akada, Toyama 939-8064, Japan

Phone: 076-420-3122 (Monday - Friday, 09:00-17:30)

Issue Date: 29th November, 2006

\_\_\_\_\_

#### Section 1: PRODUCT IDENTIFICATION

1.1. Product name: EDC-E

1.2. Product Type: Liquid Emulsion based on food-grade Vegetable Oils and

Surfactants

1.3. Hazard Rating: Health: 1 Fire: 1 Reactivity: 1

1.4. Formula: Proprietary

Substances Subject to SARA 313 Reporting Are Indicated by "#"

It is our opinion that above named product does not meet the definition of "hazardous Chemical" as defined in the OSHA "Hazard Communication Standard" regulation 29 CFR 1910.1200. This material Safety Data Sheet is provided as general information for health and safety guidelines.

#### Section 2: INGRADIENTS/IDENTITY INFORMATION

CAS No. % PEL TWA(mg/m<sup>3</sup>)

Vegetable Oil (food-grade)- proprietary Trade secret 50 15 10

Surfactants (food-grade)-proprietary Trade secret 5-10

Water 7732-18-5 45-40

#### Section 3: PHYSICAL AND CHEMICAL CHARACTERISTICS

This section completed as per formulation ingredient data unless stated

Solubility in water: dispersible

pH: 6-7

Specific Gravity: 0.94-0.96 Boiling point: 100°C for water

Melting point: NA Vapor Pressure: NA

Vapor Density: Heavier than air Percent Volatile by Volume: NA Evaporation Rate: NA

Viscosity: 20-30 cps at 20°C

Product Appearance and Odor: Off-White liquid with vegetable oil odor

.......

#### Section 4: FIRE AND EXPLOSION HAZARDS

This section completed as per formulation ingredient data unless stated.

- 4.1 Special Fire Hazards: Product-none, does not support combustion.
- 4.2 Flash point: >150°C
- 4.3 Flammable limits; Lower Explosive Limit: ND and Upper Explosive Limit: ND
- 4.4 Fire Fighting Methods: Use methods appropriate for surrounding fire.
- 4.5 Extinguishing Media: Dry chemical, CO2 or foam

Note: water, fog, and foam may cause frothing and spattering

- 4.6 Special Fire Fighting Procedures: Wear self contained breathing apparatus and Chemical resistant clothing. Use water spray to cool fire exposed containers.
- 4.7 Unusual Fire and Explosion Hazards: Burning will cause oxides of carbon

#### Section 5. HEALTH HAZARD DATA

- 5.1 This product is neither intended nor manufactured for human or animal consumption and should not be used for food or feed stuffs.
- 5.2 Effects of over exposure: NA
- 5.3 Routes of exposure: Exposure to vegetable oil mists can occur through inhalation, Ingestion, and eye or skin contact.
- 5.4 Emergency and First Aid Procedures: If inhaled, remove from contaminated atmosphere to fresh air.

Eye: immediately flush eyes with large amounts of water for 15 min. ensure rinsing entire surface of eye and under lid, if irritation persists see physician.

Lungs: move to fresh air.

Skin: wash affected areas thoroughly with soap and water.

Ingestion: Product is non-toxic, if nausea occurs, induce vomiting, and seek medical attention.

- 5.5 Effect on Humans: Most vegetable oil mists are biologically inert and are therefore considered to be nuisance particulates. Vegetable oil mists seem to have little adverse effect on the lungs and do not produce significant organic disease or toxic effects when exposures are kept under reasonable control (ACGIH 1991).
- 5.6 Carcinogenity of product-none and ingredients NTP, IARC, OSHA: No

5.7 Occupational exposure limits {8 hour time weighted average [TWA]}: mg/m3

### OSHA PEL/ACGIH TLV

Vegetable oils (Food grade)

15 (Mist)/10 (Mist)

#### Section 6. REACTIVITY DATA AND PHYSICAL HAZARDS

This section completed as per formulation ingredient data unless stated

- 6.1 Stability: Stable under normal conditions
- 6.2 Stability Conditions to Avoid: NA
- 6.3 Incompatibilities: Strong Acids and Oxidizers
- 6.4 Hazardous Decomposition Products: Product-none identified Ingredients: Carbon Oxides, Biological Decomposition (Spoilage) may result in offensive odors.
- 6.5 Hazardous Polymerization: none known
- 6.6 Polymerization Conditions to Avoid: NA
- 6.7 Vegetable oil LD 50 and LD50 Mixture: NA

.......

#### Section 7. SPILL OR LEAK PROCEDURES

This section completed as per formulation ingredient data unless stated

- 7.1 Spill Response: This product is water dispersible. Same as vegetable oil spills: isolate spill, prevent from entering water ways, and sewer systems. Sorb or remove spilled materials as soon as possible. Oils and specific quantities of oils may be reportable under government regulations.
- 7.2 Waste Disposal Method: This product is not hazardous; however, wastes must be disposed in accordance with government regulations. Consult with local sewer authority, or solid waste facility prior to disposition.
- 7.3 Precautions-Handling/Storage: Keep container closed and store cool and dry place

## **Section 8: ECOLOGICAL INFROMATION**

This section completed as per formulation ingredient data unless stated

- 8.1 Ecology: no environmental hazards are known. Avoid pollution to soil, water in accordance with government regulations
- 8.2 BOD5 AND COD: NA
- 8.3 Toxicity of the Products of Biodegradation: The product itself and its products of degradation are not toxic
- 8.4 Special Remarks on the Products of Biodegradation: NA

#### Section 9. SPECIAL PRECAUTIONS AND PROTECTION INFORMATION

No protective equipment is necessary under normal use conditions.

Precautions to be taken in handling and storage: Do not store near excessive heat or oxidizers

- 9.1 Eyes: If splashing may occur, eye protection recommended.
- 9.2 Skin: Wear impervious gloves for prolonged or repeated exposure.
- 9.3 Respiratory: Avoid breathing mists of this product

.......

#### **Section 10: TRANSPORTATION PRECAUSTIONS**

This section completed as per formulation ingredient data unless stated

Transportation considerations: This product is not classified as dangerous in the meaning of transport regulations. Shippers and transporters may need to meet packaging and transportation requirements for certain oils and respective quantities under CFR Part 130.

#### Section 11: REGULATORY INFROMATION

This section completed as per formulation ingredient data unless stated

#### 11.1 TSCA

Not listed on the TSCA inventory

Health & Safety reporting list-none of the ingredients are on the health & safety list

#### 11.2 Clean air act

This product does not contain any hazardous air pollutants This material does not contain any class 1 or 2 Ozone depleters

#### 11.3 Clean water act

None of the ingredients in this product are listed as hazardous substances, priority pollutants, or toxic pollutants under CWA

#### **Section 12: OTHER INFROMATION AND REFERENCES**

ACGIH [1991]. Documentation of the threshold limit values and biological exposure indices, 6<sup>th</sup> Ed., Cincinnati, OH, American Conference Governmental Industrial Hygienists

The above information is believed to be correct with respect to the formula used to manufacture the product in the country of origin. As data, standards, and regulations change, and conditions of use and handling are beyond our control. However, Ecocycle Corporation makes NO WARENTY, EXPRESS OR IMPLIED, IS MADE AS TO THE COMPLETENESS OR CONTINUING ACCURACY OF THIS INFORMATION. This information and product are furnished on the condition that the person receiving them shall make his/her own determination as to the suitability of the product for his/her particular purpose.