APEX COMPANIES, LLC

Remedial Investigation (RI) Report



BCP Site # C224169 2002-2024 Cropsey Avenue Brooklyn, New York

PREPARED FOR:

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Table of Contents

| 1.0 | | Intro | duction | 1 |
|-----|-----|---------|---|----|
| | 1.1 | Site D | Description | 2 |
| | 1.2 | Site S | ubsurface Geology and Hydrogeology | 3 |
| | | 2.4.1 | Regional Geology | 3 |
| | | 2.4.2 | Local Geology | 3 |
| | | 1.2.2 | Hydrogeology | 4 |
| | 1.3 | Site H | listory | 4 |
| | | 1.3.1 | Previous Investigations | 4 |
| | | | 1.3.1.1 Executive Environmental Group Phase I Environmental Site Assessment (ESA) | 4 |
| | | | 1.3.1.1 Merritt Limited Subsurface Investigation | 5 |
| | | | 1.3.1.2 Apex Limited Phase II ESA | 6 |
| 2.0 | | Reme | edial Investigation Objectives and Scope | 8 |
| | 2.1 | Reme | dial Action Objectives | 8 |
| | | 2.1.1 | Groundwater RAOs | 8 |
| | | 2.1.2 | Soil RAOs | 9 |
| | | 2.1.3 | Soil Vapor / Indoor Air Quality (IAQ) RAOs | 9 |
| 3.0 | | Reme | edial Investigation Activities & Results | 10 |
| | 3.1 | Stanc | lards, Criteria, and Guidance (SCGs) | 10 |
| | 3.2 | Qual | ity Assurance /Quality Control | 11 |
| | 3.3 | Geop | hysical Survey & Utility Markouts | 11 |
| | 3.4 | Mem | brane Interface Probe Investigation | 12 |
| | 3.5 | Soil V | apor / Indoor Air Quality (IAQ) Investigation | 12 |
| | | 3.5.1 | Upgradient Soil Vapor Points | 13 |
| | | 3.5.2 | Ambient Air Sampling | 14 |
| | 3.6 | Soil Ir | nvestigation | 14 |
| | | 3.6.1 | Dry Cleaner Area Soil Investigation | 14 |
| | | 3.6.2 | Soil Investigation During Monitoring Well Installations | |
| | 3.7 | Grou | ndwater Investigation | 16 |
| | | 3.7.1 | Initial Phase II Groundwater Investigation | |
| | | 3.7.2 | Delineation of Groundwater Impacts | |
| | 3.8 | Sumr | nary of Data Usability | 18 |
| | 3.9 | Inves | tigation-Derived Waste Disposal | 19 |

| 4.0 | | Nature and Extent of COCs and Conceptual Site Model | 20 |
|-----|-----|---|-----|
| | 4.1 | Nature and Extent of COCs in Soil | .20 |
| | 4.2 | Nature and Extent of COCs in Groundwater | .20 |
| | 4.3 | Nature and Extent of COCs in Soil Vapor and Ambient Air | .21 |
| | 4.4 | Conceptual Site Model | .22 |
| 5.0 | | Soil Vapor Mitigation Measures and IRMs | 24 |
| | 5.1 | Shopping Center SSDS IRM | .24 |
| | | Residential Building Garage Ventilation System Upgrades | |
| 6.0 | | Human Health Exposure Assessment | 25 |
| | | | |

Tables:

| Table 3-1: | Soil Vapor/Indoor Air Ambient Air Analytical Results Summary |
|------------|---|
| Table 3-2: | Soil Analytical Results Summary (VOCs) |
| Table 3-3: | Soil Analytical Results Summary (SVOCs) |
| Table 3-4: | Soil Analytical Results Summary (Pesticides, Herbicides and PCBs) |
| Table 3-5: | Soil Analytical Results Summary (Inorganics) |
| Table 3-6: | Groundwater Analytical Results Summary (VOCs) |
| Table 3-7: | Groundwater Analytical Results Summary (SVOCs) |
| Table 3-8: | Groundwater Analytical Results Summary (Pesticides, Herbicides) |
| Table 3-9: | Groundwater Analytical Results Summary (Inorganics) |

Figures:

| Figure 1: | Site Location Map |
|------------|---|
| Figure 2: | Remedial Investigation Areas |
| Figure 3: | 2002-2024 Cropsey Avenue Area Site Plan |
| Figure 4: | Well Locations |
| Figure 5: | Potentiometric Surface Maps (Includes 5A, 5B, and 5C) |
| Figure 6: | 2012 Groundwater Results Summary (Merritt) |
| Figure 7: | 2012 Soil Vapor/Indoor Air Results Summary |
| Figure 8: | 2012 Soil Sample Results Summary |
| Figure 9: | 2012 Groundwater Results Summary |
| Figure 10: | PCE Isoconcentration s in Groundwater (July 2012) |

| Figure 11: | Membrane Interface Prove Investigation Locations |
|------------|---|
| Figure 12: | Soil Vapor/Indoor Air Sample Results Summary (2012-2016) |
| Figure 13: | Supplemental Soil Investigation Results Summary |
| Figure 14: | Groundwater Investigation Results Summary (2016 Composite Data Set) |
| Figure 15: | SSDS Pilot Test Vacuum and Monitoring Point Locations |
| Figure 16: | SSDS IRM Extraction & Monitoring Points |

Appendices:

| Appendix A: | Apex Limited Phase II ESA (incorporated by reference, under separate cover) |
|-------------|---|
|-------------|---|

- Appendix B: Membrane Interface Probe Data
- Appendix C: Soil Vapor and Indoor Air Quality Analytical Reports
- Appendix D: Soil Vapor Probe Construction Information
- Appendix E: Soil Analytical Data Summary Reports
- Appendix F: Soil Boring Logs, Well Logs and Sampling Data Summaries
- Appendix G: Groundwater Analytical Data Summary Reports
- Appendix H: Data Usability Summary Report (DUSR)
- Appendix I: Garage Ventilation System Documentation.

Certification of Remedial Investigation Report

I Daniel J. Smith, P.E. of Apex Engineering, PC, certify that I am currently a New York State registered professional engineer and that this report was prepared in accordance with applicable statutes and regulations and in substantial conformance with the Division of Environmental Remediation (DER) Technical Guidance for Site Investigation and Remediation (DER-10) and that all activities were performed in accordance with the DER-approved work plan and any DER-approved modifications.



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Remedial Investigation Report 2002-2024 Cropsey Avenue Site Brooklyn, New York NYSDEC BCP# C224169

1.0 Introduction

This Remedial Investigation (RI) Report presents data related to investigation activities completed to determine the nature and extent of contaminants of concern (COCs) in soil, groundwater and soil vapor on-site at the 2002-2024 Cropsey Avenue Site (hereinafter referred to as the Site) located at 2002-2024 Cropsey Avenue, Brooklyn, New York The Site Location showing the general area of the Site is provided as **Figure 1**. The groundwater conditions associated with adjacent and nearby properties were also evaluated during the RI and summarized in this Remedial Investigation Report (RIR). The Site property and adjacent investigation areas are indicated in **Figure 2**. The primary objective of the RI is to provide an assessment of potential impacts of COCs on human health and the environment in soil, soil vapor and groundwater and to determine the appropriate remedial measures, if any, needed for the Site.

The application to the NYSDEC Brownfield Cleanup Program (BCP) was submitted on behalf of 2002 Cropsey Associates, LLC (the Participant) and was accepted on August 4, 2014. The Brownfield Cleanup Agreement (BCA) was executed with the NYSDEC on September 30, 2014. The RI was completed in general accordance with the Remedial Investigation Work Plan (RIWP) approved by the New York State Department of Environmental Conservation (NYSDEC).

Environmental Site Characterization (ESC) investigations were completed at the Site by multiple parties between 2004 and 2013. Based on the findings of the ESCs and a June 2, 2014, meeting between the NYSDEC, the Participant and Apex, RI activities were conducted on the following three parcels:

- 2002 2024 Cropsey Avenue (Block 6467, Lot 1). This property is the Site which is currently occupied by a small strip mall / shopping center fronting Cropsey Avenue. The shopping center includes a former and current dry-cleaning establishment with partial basement and several small exterior access ways for shop owners at the rear of the site.
- 8831 and 8841 20th Avenue (Block 6467, Lot 12). This property includes residential buildings located to the southwest of the Site. The Block 6467, Lot 12 property is separated from the shopping center structure by a small grassy area between the strip mall building structure and the residential building towers. The residential towers contain sub-grade parking facilities.
- 2036 Cropsey Avenue (Block 6469, Lot 1). This property contains residential buildings to the southeast of the Site across Bay 25th Street. The general construction is similar to the Block 6467, Lot 12 properties with a subgrade parking garage and overlying residential towers.

The general areas investigated were located along Cropsey Avenue, 20th Avenue and Bay 25th Street (see Figure 2).

The results of the ESCs and the initial remedial investigations identified the presence of volatile organic compounds (VOCs) in soil vapor underlying the strip mall building at the Site that exceeded the New York State Department of Health (NYSDOH) Matrix 1 and Matrix 2 thresholds for Sub-Slab Vapor Intrusion Mitigation as required by the NYSDOH

"Guidance for Evaluating Soil Vapor Intrusion in the State of New York", October 2006 (as amended) (herein referred to as the NYSDOH Guidance Document).

Several mitigation measures were implemented (and are being implemented) to control soil vapors from potentially entering the occupied spaces of the Site buildings and the residential towers on the adjacent properties. The parking garage systems underlying the Block 6467, Lot 12 and Block 6469, Lot 1 residential towers were upgraded with improved air-moving equipment / ventilation systems in July - August 2016 to mitigate potential vapor intrusion into overlying residential areas. In addition, a NYSDEC-approved Interim Remedial Measure (IRM) was proposed and approved by the NYSDEC on January 27, 2017, to install a Sub-Slab Depressurization System (SSDS) to mitigate soil vapor intrusion into the shopping center located at the Site. This shopping center IRM is to be installed in September 2017, following procurement and approvals needed to proceed.

The following sections of the RIR present data identifying the nature and extent of soil, soil vapor and groundwater impacts at the site and the immediately adjacent properties within the approved scope of work and the basis for future remedial actions.

1.1 Site Description

The Site is located at 2002-2024 Cropsey Avenue in the Borough of Brooklyn, New York City, New York and occupies a parcel that is identified by Tax Map Number: Block 6467, Lot 1. (see **Figures 1, 2 and 3**). The Site is currently improved with a single-story, multi-unit retail building (which has a partial basement) with an approximately 15,000 square foot (ft²) area. Concrete-paved sidewalks surround the Site building to the northwest, northeast and southeast. There is a small parking area / trash enclosure located at the southwest corner of the building. There is also a small rear access way to the dry-cleaning establishment within the shopping center at the southeast corner of the building (see **Figure 3**). The southwest exterior building wall defines the boundary of the BCP Site as indicated in **Figure 3**.

The Site is bounded by Cropsey Avenue to the northeast, 20th Avenue to the northwest, a residential building with subgrade parking to the south and southwest, and Bay 25th Street to the southeast. The elevation of the Site is approximately 20 feet above mean sea level (msl). Surface topography consists of a gentle downward slope to the southwest towards Gravesend Bay, which is approximately 1,000 feet from the Site. A narrow undeveloped strip of land extends along the entire south (rear) side of the Site building and is not part of the BCP Site. The layout of the Site and surrounding properties is presented in **Figure 2** and **Figure 3**. Local groundwater flow in the vicinity of the Site is generally to the south with localized southeast and southwest components towards Gravesend Bay. The Site-specific depth to water is variable but generally is approximately 15 and 20 feet below grade surface (bgs). Additional information regarding local soil and groundwater properties is provided in **Section 1.2**.

Currently, the Site is developed with a strip-retail shopping center. The Site, and the adjacent residential properties included in this RI, are located within an R6, Residential Zone district. Land use at the Site is commercial and land use at the adjacent properties include din this RI is residential. Existing shopping center tenants include a luncheonette and restaurant, a nail salon, convenience / drug store, distributors / traders, and a dry cleaner (not the same as the historic dry cleaner associated with contamination at the site). The current dry cleaner operation, GLY Cleaners, consists of a closed loop hydrocarbon cleaning system. GLY Cleaners also offers tailoring services. Historic dry-cleaning operations which utilized tetrachloroethene (PCE) were conducted in the existing tenant space and in the partial basement of the Site.

1.2 Site Subsurface Geology and Hydrogeology

This section of the RIR describes the regional / local geology and hydrogeology.

2.4.1 Regional Geology

The unconsolidated geologic deposits underlying Brooklyn consist of clay, silt, sand, and gravel that overlie southward-dipping consolidated bedrock. The crystalline bedrock consists mainly of Precambrian age granite, gneiss, and schist. The overlying unconsolidated sediments were deposited during the Cretaceous and form, in a descending order, the Raritan and Magothy Formations.

During the Pleistocene, several episodes of glaciation eroded the Cretaceous deposits (Smolensky, et al, 1989). The oldest Pleistocene deposit is the Jameco Gravel (Jameco aquifer), which overlies the Magothy Formation and Raritan Clay and is present only in western Long Island. The Gardiners Clay overlies the Jameco Gravel, Magothy Formation, and Raritan Clay (a confining unit). The Upper Pleistocene deposits formed when the glacial ice and glacial melt water deposited till and outwash material, forming what is presently known as the Upper Glacial aquifer. The Raritan Formation consists of the Lloyd Sand and the Raritan Clay. The Lloyd aquifer (the hydrogeologic equivalent of the Lloyd Sand) consists of fine to coarse sand, gravel, commonly with a clayey matrix, and lenses and layers of silty and solid clay. The Raritan Clay is regionally continuous and consists of silty and solid clay, and lenses and layers of sand. Because of its low permeability, the Raritan Clay serves as a confining unit for the underlying Lloyd Sand.

The Magothy Formation is a deltaic deposit consisting of fine to medium sand, clayey in part, interbedded with lenses and layers of coarse sand, silt, and sandy and solid clay. Gravel is common in the basal zone of the Magothy Formation. The Jameco aquifer (the hydrogeologic equivalent of the Jameco Gravel) is a channel filling consisting of fine to very coarse sand and gravel with few layers of clay and silt (Smolensky, et al, 1989). The Gardiners Clay is a lagoonal/shallow-bay clay consisting of clay, silt, and few layers of sand and gravel (Smolensky, et al, 1989). The Upper Glacial aquifer consists primarily of till and glacial outwash deposits. The till, composed of clay, sand, gravel, and boulders, forms the Harbor Hill and Ronkonkoma terminal moraines. These terminal moraines represent the farthest advance of late-Pleistocene glaciation on Long Island. South of the moraine deposits is a glacial outwash plain, which, in Kings County, extends from the Harbor Hill moraine to Jamaica Bay and New York Bay, and consists of fine to very coarse sand and pebble to boulder sized gravel (Smolensky, et al, 1989).

2.4.2 Local Geology

The Site is located south of the Harbor Hill terminal moraine and the surficial deposits consist of glacial outwash deposits (Upper Glacial aquifer) at the Site. Based on a review of the U.S. Geological Survey publication titled Hydrologic Framework of Long Island, New York, U.S. Geological Survey Hydrologic Investigations Atlas HA-709 (Smolensky, et al, 1989), bedrock beneath the Site is expected to occur approximate elevation of 650 feet below msl. The Lloyd aquifer, which overlies bedrock, has a surface elevation of approximately 500 feet below msl. The Raritan Clay has a surface elevation of approximately 400 feet below msl. The Magothy aquifer has a surface elevation of approximately 250 feet below msl. The Jameco aquifer has a surface elevation of approximately 200 feet below msl. The Gardiners Clay has a surface elevation of approximately 150 feet below msl. The Upper Glacial aquifer corresponds to the saturated upper part of the highly permeable Pleistocene deposits of sand and gravel.

Based on the soil borings installed during the ESCs and RI, fine to coarse sand deposits (glacial outwash deposits [Upper Glacial aquifer]) were encountered. Apex encountered approximately five (5) to seven (7) feet of fill in the upper portion of the soil borings. No confining layers were observed during the drilling activities

1.2.2 Hydrogeology

The principal aquifers underlying the Site are the Upper Glacial aquifer, Jameco aquifer, and Magothy aquifer. The Gardiners Clay hydraulically confines the Magothy and Jameco aquifers in most of Brooklyn; the Jameco aquifer and Magothy aquifer hydrogeologic units are in direct hydraulic connection with each other.

Groundwater in the Upper Glacial aquifer occurs under unconfined conditions at and near the Site. Within the project area, the average horizontal hydraulic conductivity of the Upper Glacial aquifer is approximately 270 feet per day (ft./d), with an anisotropy ratio of approximately 10:1 (horizontal to vertical, respectively) (McClymonds and Franke, 1972). The average horizontal hydraulic conductivity of the Jameco aquifer in the project area is approximately 200 to 300 ft./d, with an anisotropy ratio of approximately 10:1 (horizontal to vertical, respectively) (McClymonds and Franke, 1972). The average horizontal hydraulic conductivity of the Magothy aquifer in the project area is approximately 50 ft./d, with an anisotropy ratio of approximately 100:1 (horizontal to vertical, respectively) (McClymonds and Franke, 1972). The average horizontal hydraulic conductivity of the Magothy aquifer in the project area is approximately 50 ft./d, with an anisotropy ratio of approximately 100:1 (horizontal to vertical, respectively) (McClymonds and Franke, 1972).

The Site is located approximately 1,000 feet northeast of Gravesend Bay. Based on data collected from monitoring wells installed during the RI, groundwater occurred at an average depth of 20-feet bgs and the shallow groundwater flow was to the south / southeast (see Figure 4 for monitoring well locations and Figure 5 (including Figures 5A, 5B, and 5C) for potentiometric surface maps). In the intermediate zone, a more southwesterly flow component appeared to be evident based upon the data available.

1.3 Site History

Based on review of available historical information, the Site was vacant land prior to the construction of the current on-site structures c. 1950 and the on-site building structure has been relatively unchanged since 1950. At least four dry cleaners have operated at the Site, including the following:

- Augie's Cleaners (1991 to1996);
- Michael's Cleaners (1996 to 2005);
- Ida Cleaners (2005 to 2007); and,
- GLY Cleaners (2007 to current).

No information is available regarding tenants at the Site prior to 1991. Based upon available records, it appears that all historic dry-cleaning operations were conducted in the same physical tenant space, hereinafter referred to as the Dry Cleaner Area.

1.3.1 Previous Investigations

The following sections provide a summary of the ESC activities conducted for the Site to date, including those conducted by Merritt Environmental Consulting Corp. (Merritt) circa 2012 and by Apex. Apex does not make any comment regarding the accuracy of work performed by others and is reporting historical investigation results as a basis for the work later completed by Apex on behalf of the Participant. Apex is not responsible for any inaccuracies in the reporting done by others.

1.3.1.1 Executive Environmental Group Phase I Environmental Site Assessment (ESA)

Executive Environmental Group (EEG) completed an Environmental Site Assessment - Phase I of the Site on March 10, 2004. EEG identified a commercial site consisting of eight stores and a total of 14,691 square feet. This assessment

was conducted by Eliot Ely, CEI, who is an EPA certified asbestos inspector, EPA certified lead based paint supervisor and certified by the Environmental Assessment Association as an Environmental Inspector.

The scope of this Site Assessment included a visual inspection, governmental database review, Sanborn Map review, interviews and related sources. The purpose of this inspection was to review the range of common hazardous contaminants and to evaluate recognized environmental conditions at the subject site as per ASTM Practice E-1527-00. EEGs professional opinion regarding the retail space and its basement did not show any environmental concerns and concluded the following:

- Past and Current Uses of Premises: The Site was being exclusively used as a commercial building since 1950. Prior to this the site was a vacant piece of land. There was no historical evidence or known evidence of environmental hazards associated with the past or current tenant use of this subject site reported by EEG.
- Asbestos Containing Materials (ACM) and Lead based Paint (LBP): A survey for friable ACM and LBP was
 conducted at the site in conjunction with the Phase I survey. There are no ACM or LBP concerns noted at the
 subject site by EEG.
- Storage Tanks: There were no aboveground storage tanks (ASTs) or underground storage tanks (USTs) noted at the time of EEG's site visit and Phase I ESA.
- Hazardous Materials Use: EEG identified an operating dry cleaner in 2004 (Michael's Drycleaners, USEPA ID No. NYD093768539). EEG indicated there were no violations associated with this operation on the database sources it reviewed. Manifests showing proper disposal of chemicals generated via Safety Kleen were noted by EEG. No evidence of spills or staining were noted by EEG at the time of their site visit.
- General EEG Conclusion: "There were no environmental concerns found at the site."

1.3.1.1 Merritt Limited Subsurface Investigation

Merritt conducted a limited Subsurface Investigation in February 2012 to evaluate groundwater quality conditions downgradient of the Dry Cleaner Area (see **Appendix A** by reference under separate cover). The Merritt report indicated that the groundwater flow direction in the vicinity of the Site was not determined during their investigation, but was believed to be "from roughly north to south towards Gravesend Bay" which would place the current GLY Cleaners facility (and location of former dry cleaning operations that were the suspected source of environmental impact) hydraulically upgradient of the adjacent residential properties to the south (i.e., 8831 and 8841 20th Avenue property and 2036 Cropsey Avenue site). It appears that two of the three temporary monitoring points installed by Merritt (i.e., B2 (Merritt) and B3 (Merritt)) were completed at the side yard of the 8831 and 8841 20th Avenue property in what was believed by Merritt to be hydraulically downgradient of the Dry Cleaner Area.

The soil samples collected as part of well installation were field screened by visual / olfactory inspections and for the presence of total VOCs utilizing a photo-ionization detector (PID). Merritt indicated that none of the soil samples exhibited field-screening evidence of VOC-related impacts; however, none of the soil samples were submitted for laboratory analyses. Tetrachloroethylene (PCE) was detected in all three groundwater samples with the highest concentrations in the groundwater samples collected from B2 (Merritt) and B3 (Merritt) at 720 micrograms per liter (µg/L) and 63,000 µg/L, respectively. The B3 (Merritt) groundwater sample was collected from the rear of the TSOC in an estimated downgradient location. The NYSDEC Class GA Groundwater Quality Standard (GWQS) for PCE is 5 µg/L. Sample results from the Merritt investigation are summarized in **Figure 6**.

Following review of data generated by Merritt, which indicated elevated concentrations of COCs in the groundwater samples, the NYSDEC was notified and Spill No. 11-13648 was assigned to the Site.

1.3.1.2 Apex Limited Phase II ESA

As part of a Limited Phase II ESA to better define the investigation results reported by Merritt, Apex completed soil vapor intrusion (SVI), soil and groundwater investigations on the Site and portions of the immediately adjacent property in July 2012. Apex installed and sampled three interior soil borings, six exterior soil borings, three groundwater monitoring wells and five sub-slab soil vapor probes. The Apex Limited Phase II ESA Report is included under separate cover by reference as **Appendix B** and is summarized below.

SVI Investigation

Sub-slab soil vapor (SV) and indoor air quality (IAQ) samples were collected from five co-located locations within the Dry Cleaner Area (current GLY Cleaners and same location as former dry-cleaning operations) and at the residential parking garage located closest to the TSOC (8841 20thAvenue Site) to the south-southwest (downgradient) of the Site. The nearby building on the 8841 20thAvenue Site was a multi-level residential apartment building with a subterranean, basement parking garage. No residential units were observed in the basement area which is utilized for parking and tenant-storages purposes.

Two sampling points were located within the Dry Cleaner Area. Per NYSDOH protocols, collocated SV and IAQ samples were collected at each location. As shown on **Figure 7**, IAQ-1 / SV-1 and IAQ / SV-2 were collected from within the basement of the Dry Cleaner Area on the Site. Similarly, three sets of samples (IAQ and SV) were collected at the adjacent residential parking garage (i.e., IAQ-3/SV-3, IAQ-4/SV-4 and IAQ-5/SV-5). **Figure 7** presents the locations of all the sampling points.

The SV / IAQ samples were collected in six-liter Summa canisters over an eight-hour period and analyzed by a NYSDOH Environmental Laboratory Accreditation Program (ELAP)-certified laboratory for NYSDEC Target Compound List (TCL) VOCs by EPA Method 8260. The samples were analyzed in accordance with Level A Analytical Services Protocols (ASP).

The SVI sampling results indicated the presence of elevated PCE and trichloroethene (TCE) concentrations in the soil vapor underlying the Dry Cleaner Area and the parking garage. The highest SV concentrations of PCE and TCE detected were 420,000 micrograms per cubic meter (μ g/m³) and 6,600 μ g/m³, respectively in the SV-2 sampling location located in the basement of the Dry Cleaner Area. The highest sub-slab, soil vapor concentrations of PCE (210,000 μ g/m³) and TCE (790 μ g/m³) in the parking garage area were detected in the samples collected from SV-4 located within the basement of the 8841 20th Avenue Site.

The highest concentrations of COCs in the indoor air samples were from Sample IAQ-2 in the basement of the Dry Cleaner Area (PCE was detected at 100 μ g/m³ and TCE was detected 18 μ g/m³). The IAQ samples collected within the parking garage at the 8841 20th Avenue Site did not exhibit elevated concentrations of PCE or TCE. The maximum concentrations of the parking garage IAQ samples were 4.3 μ g/m³ of PCE and 1.1 μ g/m³ of TCE. (see **Figure 7** and **Appendix A**).

Based on the concentrations of COCs detected in the sub-slab soil vapor samples, further delineation in both soil vapor and groundwater matrices to the south and east was recommended. The additional delineation of the extent of elevated soil vapors of PCE and TCE is discussed later in this RIR.

Soil Sampling

Twelve borings, with associated soil sampling and analyses, were advanced between July 11 and July 13, 2012, (see

Figure 8). As indicated in **Figure 8**, three soil borings were conducted within the basement of the Dry Cleaner Area and the remaining borings were conducted in the soil / grass -covered area just to the southwest of the Site on the adjacent 8841 20thAvenue Site. Based upon field-screening results, one sample per boring was submitted to the laboratory and analyzed for TCL VOCs by EPA Method 8260. For the purposes of this ESC, the soil analytical data were compared to the NYSDEC Unrestricted Use Soil Cleanup Objectives (SCOs) included in Part 375-5-6.8 - NYSDEC Remedial Program Soil Cleanup Objectives.

In general, only low concentrations of PCE and TCE were detected in the associated soil samples, with the exception of the S-1 S (i.e., "shallow") sampling located in the basement of the Dry Cleaner Area. PCE was detected at 14,000 micrograms per kilogram (μ g/kg) at a depth of 3.0 to 4.0 feet bgs in the S-1 S sample, which exceeded the NYSDEC Unrestricted Use SCO for PCE of 1,300 μ g/kg. All other TCL VOCs in the remaining 11 soil sample locations were either: not detected at concentrations above their respective laboratory method detection limits (MDLs); detected at concentrations in laboratory blank samples as well as the characterization samples indicating that they were laboratory artifacts.

Groundwater Sampling

Apex collected nine samples on July 13, 2012 to assess groundwater quality conditions; three from the newly installed monitoring wells (MW-1, MW-2 and MW-3) and six from temporary wells that originated as soil borings. All wells during this sampling event were screened in the shallow groundwater zone, within approximately 5 to 10 feet of the top of the water table. The groundwater samples were submitted to the NYSDOH ELAP-certified laboratory for NYSDEC and analyzed for TCL VOCs by EPA Method 8260 and TCL semi-volatile organic carbons (SVOCs) by EPA Method 8270 in accordance with Level A ASP.

For the purposes of this ESC, the groundwater analytical data were compared to the NYSDEC Class GA Groundwater Quality Standards (GWQS) and Guidance Values set forth in the NYSDEC Division of Water Technical and Operational Guidance (TOGS) Series 1.1.1 Ambient Water Quality Standards (AWQS) and Guidance Values and Groundwater Effluent Limitations – Reissued June 1998 and April 2000 Addendum.

As summarized in **Figure 9**, PCE was detected in all groundwater samples at concentrations exceeding its Class GA GWQS of 5 micrograms per liter (µg/L). The highest concentrations of PCE were detected in the groundwater samples collected from monitoring wells MW-1 (2,500 µg/L) and MW-2 (1,400 µg/L), which are both located directly downgradient of the Dry Cleaner Area. TCE and cis-1,2-dichlorethene (a PCE degradation product) were also both detected at concentrations above their respective Class GA GWQSs in the two aforementioned groundwater samples.

A PCE isoconcentration map was prepared (**Figure 10**) to illustrate that the highest concentration of TCL VOCs was detected in the groundwater samples collected from hydraulically downgradient of the Dry Cleaner Area on the bordering 8841 20thAvenue Site. PCE (concentrations exceeded 1,000 μ g/L) and cis-1,2-DCE concentrations at 1,300 μ g/L) were both detected in the groundwater sampling locations located along the southeast property line indicating the probable off-site migration of a COC groundwater plume to the south and east of the Dry Cleaner Area, consistent with the inferred groundwater flow direction.

The groundwater sample collected from the upgradient monitoring well MW-3 contained PCE at a concentration of 14 μ g/L indicating the presence of a minor source of contaminants upgradient of the Site. That is likely contributing to the overall distribution of PCE and its degradation products in the investigation area. All TCL SVOCs were detected at concentrations below their respective laboratory MDLs or below their respective Class GA GWQ standards or guidance values.

2.0 Remedial Investigation Objectives and Scope

As summarized above, the ESCs completed for the Site indicated the presence of COCs in exceedance of NYSDEC applicable standards, criteria and / or guidelines in soil, groundwater, soil vapor and indoor air. Therefore, per NYSDEC Department of Environmental Remediation (DER) protocols, the conduct of an RI was required to evaluate the nature and extent contamination in the soil, soil vapor and groundwater matrices both on- and off- Site. The subsequent RI was completed at this Site in accordance with the July 2015 RIWP approved by the NYSDEC.

The objectives of this RI were to:

- Determine the nature and extent of COCs in soil, groundwater and soil vapor on Site, as well as on nearby properties hydrogeologically downgradient of the Site where preliminary data indicated off-site impacts may be present and related to the Site impacts previously identified;
- Develop a further understanding of the local groundwater conditions including potential fate and transport of COCs emanating from the Dry Cleaner Area;
- Assess potential impacts to human health (qualitatively) and the environment (qualitatively and quantitatively) as a result of the release of COCs at the Site; and,
- Collect and process data sufficient to support making decisions regarding remediation of the Site and to serve as a design basis for future evaluation of potential remedial actions.

2.1 Remedial Action Objectives

Remedial Action Objectives (RAOs) were established prior to commencing the RI and were presented in the July 2015 RIWP to ensure that data obtained during the RI were directly related to the defining future remediation needs at the Site, if necessary. In accordance with the NYSDEC DER-10, RAOs considered the following:

- Applicable Standards, Criteria and Guidance (SCGs) which considered the current, intended and reasonably anticipated future use of the Site and its surroundings;
- All contaminants which were detected at concentrations exceeding applicable SCGs;
- The types of environmental media impacted by such contaminants;
- The extent of the impact to the environmental media;
- Actual or potential human exposures and / or environmental impacts resulting from the contaminants in environmental media identified above; and,
- Any site-specific cleanup levels which were developed.

The RAOs identified in Section 2.1.1 through Section 2.1.3 are applicable at the Ste and are incorporated into the RIR.

2.1.1 Groundwater RAOs

The following RAOs shall apply to groundwater:

- RAO for Public Health Protection: (1) Prevent / mitigate contact with, or inhalation of, volatiles from contaminated groundwater; and,
- RAO for Environmental Protection: (1) Restore the groundwater aquifer to predisposal / pre-release conditions or applicable regulatory criteria to the extent practicable; and, (2) Remove the source of ground impacts to the extent practicable.

2.1.2 Soil RAOs

The following RAOs shall apply to soil:

- RAO for Public Health Protection: (1) Prevent / mitigate ingestion and / or direct contact with impacted soil; and, (2) Prevent / mitigate inhalation of, or exposure from, COCs volatilizing from contaminants in soil.
- RAO for Environmental Protection: (1) Prevent migration of COCs that would result in groundwater or surface water contamination.

2.1.3 Soil Vapor / Indoor Air Quality (IAQ) RAOs

The following RAOs shall apply to soil vapor:

• *RAO for Public Health Protection:* (1) Mitigate impacts to public health resulting from existing, or the potential for, soil vapor intrusion into the building(s) at the site and adjacent properties.

3.0 Remedial Investigation Activities & Results

Based on the detection of COCs in the sub-slab soil vapor, indoor air, soil and groundwater samples collected during the ESCs, further investigation was required to evaluate the nature and extent of the environmental impacts. The RI activities discussed in this RIR were completed per the RIWP and/or communications with the NYSDEC and in general accordance with the Field Sampling Plan (FSP), Quality Assurance Project Plan (QAPP) and Health and Safety Plan (HASP) that were incorporated within the NYSDEC-approved RIWP.

3.1 Standards, Criteria, and Guidance (SCGs)

SCGs are defined as "mean standards and criteria that are generally applicable, consistently applied, and officially promulgated, that are either directly applicable, or that are not directly applicable but are relevant and appropriate, unless good cause exists why conformity should be dispensed with, and with consideration being given to guidance determined, after the exercise of scientific and engineering judgment, to be applicable."

The most common SCGs applicable in New York State, as promulgated by the NYSDEC and / or NYSDOH and at the Site include the following:

- **Soil**: SCOs and supplemental SCOs identified in 6 NYCRR 375-6.8 and the Commissioner's Policy on Soil Cleanup Guidance (CP- 51);
- **Groundwater**: Class GA Groundwater Quality Standards and Guidance Values set forth in the NYSDEC Division of Water Technical and Operational Guidance Series 1.1.1 Ambient Water Quality Standards and Guidance Values and Groundwater Effluent Limitations Reissued June 1998; and,
- Soil Vapor / Indoor Air Quality: NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York, October 2006, as amended including updates in September of 2013 and August of 2015.

In addition to the regulatory SCGs identified above, the follow regulations also apply to the remedy selection and implementation process:

- New York Codes, Rules and Regulations (NYCRR) Part 175 -Special Licenses and Permits --Definitions and Uniform Procedures;
- NYCRR Part 371 -Identification and Listing of Hazardous Wastes;
- NYCRR Part 372 Hazardous Waste Manifest System and Related Standards for Generators, Transporters and Facilities (November 1998);
- NYCRR Subpart 374-1 Standards for the Management of Specific Hazardous Wastes and Specific Types of Hazardous Waste Management Facilities;
- NYCRR Part 375 Environmental Remediation Programs;
- NYCRR Part 376 Land Disposal Restrictions;
- NYCRR Part 608 Use and Protection of Waters;
- NYCRR Parts 700-706 Water Quality Standards;
- NYCRR Part 750 through 758 Implementation of NPDES Program in NYS (SPDES Regulations);
- Code of Federal Regulations (CFR) Part 1910.120 Hazardous Waste Operations and Emergency Response;
 and CFR Part 144 Underground Injection Control Program;

Formal regulations are not the only SCGs that may be applicable. The following regulatory guidance documents may be applicable and will be considered in the final remedy design and implementation:

- United States Environmental Protection Agency (USEPA) Office of Solid Waste and Emergency Response (OSWER) Directive 9355.047FS Presumptive Remedies: Policy and Procedures;
- USEPA OSWER Directive 9355.048FS Presumptive Remedies: Site Characterization
- and Technology Selection for CERCLA sites with Volatile Organic Compounds in Soils;
- DER -10 Technical Guidance for Site Investigation and Remediation;
- DER-15 Presumptive/Proven Remedial Technologies;
- Technical and Administrative Guidance Memorandum (TAGM) 4013 Emergency Hazardous Waste Drum Removal/ Surficial Cleanup Procedures;
- TAGM 4059 Making Changes to Selected Remedies;
- TAGM 3028 "Contained In" Criteria for Environmental Media: Soil Action Levels;
- TOGS 1.1.1 Ambient Water Quality Standards & Guidance Values and Groundwater Effluent Limitations;
- TOGS 1.3.8 New Discharges to Publicly Owned Treatment Works;
- TOGS 2.1.2 Underground Injection/Recirculation (UIR) at Groundwater Remediation Sites;
- Commissioners Policy (CP) 43 Groundwater Monitoring Well Decommissioning Procedures;
- CP-51 Soil Cleanup Guidance;
- Air Guide 1 Guidelines for the Control of Toxic Ambient Air Contaminants;
- Citizen Participation in New York's Hazardous Waste Site Remediation Program: A Guidebook;
- OSWER Directive 9200.4-17 Use of Monitored Natural Attenuation at Superfund, RCRA Corrective Action, and Underground Storage Tank Sites; and,
- NYSDOH Environmental Health Manual CSFP-530 "Individual Water Supplies Activated Carbon Treatment Systems."
- NYSDOH Guidance for Evaluating Soil Vapor Intrusion in the State of New York (as amended)

It should be noted that these SCGs are only potentially applicable at this stage of the RI program and a more detailed applicability review will be completed as part of the future Remedial Design / Remedial Action (RD/RA) process.

3.2 Quality Assurance /Quality Control

A QAPP was prepared and submitted with the RIWP and was approved by the NYSDEC. The QAPP was prepared in accordance with the USEPA guidance entitled Guidance for Quality Assurance Project Plans EPA QA/G-5 (USEPA, 2002), the NYSDEC DER-10 Technical Guidance for Site Investigation and Remediation (NYSDEC, 2010), and considering requirements of the NYSDEC BCP program.

The objective of the QAPP was to ensure that the data produced as a result of the various sampling and monitoring activities, including soil, groundwater, soil vapor, and ambient (indoor and outdoor) air matrices was of the highest quality and usable for the intended purpose of making decisions related to the Site.

3.3 Geophysical Survey & Utility Markouts

Prior to commencing any subsurface remedial investigation activities, a Site and utility inspection was completed;

and the proposed sampling and/or monitoring well locations were marked accordingly in the field. After the locations were pre-selected, New York's Dig Net of New York City & Long Island was contacted to mark underground utilities and a subcontractor was retained to conduct a private utility mark-out and clear each subsurface location for the presence of sub-grade utilities. Site property owners and adjacent property owners were contacted; as-builts drawings were reviewed and personnel with knowledge of subsurface utilities and structures were contacted. The borehole locations were also assessed using ground penetrating radar (GPR) to ensure each location was clear of utilities or subsurface structures. Utility mark-outs were completed on September 23, 2015, March 7, 2016 and June 22, 2016. There were no utility contacts or issues related to health and safety during performance of the RI.

The following sections provides a detailed discussion of the investigation work-flow components conducted as part of the RI. Summaries of the resultant analytical data are also presents; however, more detailed discussions of the natural and extent of contamination (i.e., interpretation of factual data) are presented in **Section 4**.

3.4 Membrane Interface Probe Investigation

As discussed in the NYSDEC approved RIWP, a Membrane Interface Probe (MIP) investigation was conducted on October 27 and 28, 2015 to assess the distribution of chlorinated volatile organic carbons (CVOCs) in the subsurface unsaturated and saturated soils. The MIP data were to be utilized to confirm the final monitoring well screened interval depths. The MIP locations are detailed on **Figure 11**.

The MIP system included a PID and flame ionization detector (FID) for the detection of VOCs and a halogen-specific electron detector (XSD) for the detection of CVOCs. As the probe was advanced into the subsurface, it continuously measured soil conductivity using the XSD and VOCs using the PID and FID. The probes were advanced to nominal depths of approximately 40 feet bgs. The data from the October 2015 MIP investigation is presented in **Appendix C**.

As reported in the December 3, 2015 correspondence with the NYSDEC, the results of the MIP investigation did not indicate elevated concentrations of CVOCs in the areas assessed, even near the MW-1 location, where elevated concentrations of CVOCs were previously detected in groundwater samples. Therefore, a modified monitoring well installation plan was approved by the NYSDEC in a December 18, 2015 telephone conversation with Apex. Using a phased in approach, two intermediate depth groundwater monitoring wells were installed at the previously-identified source area located near monitoring well MW-1 and upgradient of the source area located near MW-4. A full round of synoptic groundwater elevation data and COC analytical data was then used to determine the final delineation locations (see **Figure 5**) for the site-specific potentiometric surface maps prepared as part of this task.

Groundwater delineation activities are discussed further in Section 3.6 below.

3.5 Soil Vapor / Indoor Air Quality (IAQ) Investigation

As noted above, previous ESCs reported the presence of PCE and TCE at elevated concentrations in both sub-slab SV samples and IAQ samples at the Dry Cleaner Area, the Site and the adjacent parking garage. Therefore, based on the results of the 2012 ESCs, further characterization and delineation of the soil vapor and indoor air for both the Site and the adjacent and nearby properties were conducted, as discussed below.

Figure 12 summarizes the sub-slab SV and IAQ investigations completed between 2012 (ESC data) and 2017 (RI data). RI soil vapor points SV-9 through SV-11 were installed on October 26, 2015, and samples were collected on October 29, 2015. Sub-slab and soil vapor points SV-6 through SV-8 were installed on October 29 and samples were collected on October 30, 2015. In addition, soil vapor locations SV-3, SV-4, and SV-5 were re-sampled as part of the RI on April 14, 2017, to assist in evaluating effectiveness of the upgraded residential parking garage ventilation fan systems. The soil vapor probes were installed as summarized below. Results of sampling events are provided in **Table 3-1** and **Figure 12**.

Shopping Center Vapor Probes:

- SV-1, SV-2, and SV-6 (interior sub-slab probes with collocated indoor air samples [IAQ-#] at SV-1 and SV-2);
- SV-7, SV-8, and SV-12 (exterior probes on the northwest, north, and northeast sides of the building). SV-12 also serves as an upgradient probe relative to the Dry Cleaner Area.

8831 and 8841 20th Avenue Vapor Probes: *

• SV-3, SV-4, SV-5, SV-9, and SV-10 (interior sub-slab probes with collocated indoor air samples);

2036 Cropsey Avenue Vapor Probes: *

• SV-11 (interior sub-slab probe with collocated indoor air sample);

Off-site, Upgradient Background Vapor Probes:

SV-13 and SV-14 (exterior sub-slab probes).

* It should be noted that the ventilation fan upgrades in the residential parking garage were completed on or before October 8, 2015. All vapor samples prior to that date were collected without the residential parking garage fans running and all samples collected after that date were collected with the residential parking garage fans running under normal operating conditions.

In accordance with the NYSDEC-approved QAPP, the soil vapor, indoor air and ambient air samples, and associated QA/QC samples were collected in six-liter Summa canisters over an eight-hour period and submitted to SGS Accutest Laboratories (SGS), a NYSDOH ELAP-certified laboratory for CVOC analyses by EPA Method TO-15. The original laboratory data sheets are included in **Appendix C**¹.

For the purposes of this RIR, the sub-slab soil vapor data, indoor air quality data and ambient air data are compared to the thresholds included in Matrices 1 and 2 included in the NYSDOH Guidance Document (as amended and applicable). The indoor air concentrations are also compared to the NYSDEC Air Guideline Values for PCE of 30 μ g/m³ and TCE of 2 μ g /m³. Results of the soil vapor and indoor air sample analysis, as well as the appropriate SCGs, are summarized on **Table 3-1** and **Figure 12**. Additional information regarding sample results, including the most recent April 2017 data, is provided in **Section 4.3**.

3.5.1 Upgradient Soil Vapor Points

Apex installed soil vapor points SV-12 through SV-14 in March 2016 to assess soil vapor conditions outside of building footprints in locations hydrogeologically upgradient of the Site. The three upgradient soil vapor points were installed to depths of 5.0 feet below grade surface (bgs) with six-inch screens. The annular spaces surrounding the wells

¹ Lab analytical summary reports only are included in the Appendices for all media. Complete analytical packages for all media sampled are provided electronically and are discussed further in the Data Usability Summary Report.

screens were backfilled with clean filter-pack sand, followed by a bentonite seal and cement grout (see wellconstruction diagrams in **Appendix D**. Soil vapor and associated QA / QC samples were collected on March 17, 2016, per the NYSDEC-approved QAPP. As summarized in **Table 3-1** the following conditions were noted upgradient of the Site.:

- The SV-12 soil vapor sample contained vinyl chloride (VC) at a con centration of 854 μg/m³;
- The SV-13 soil vapor sample contained only low-level concentrations of the chlorinated VOCs of concern (14 μg/m³ PCE and non-detectable levels of TCE, DCE and vinyl chloride); and,
- The SV-14 soil vapor sample exhibited PCE at concentrations at 258 µg/m³.

3.5.2 Ambient Air Sampling

Ambient, background air samples were collected on March 9, 2016, and March 17, 2016, and analyzed for VOCs by EPA Method TO-15. Six-liter summa canisters collected air over an eight-hour period. One sample was collected from a hydraulically downgradient of the Dry Cleaner Area near monitoring well MW-11 on March 9, 2016. The March 17th sample was collected from a hydraulically upgradient location in the vicinity SV-14. Very low concentrations of COCs were detected in the March 9th sample (all chlorinated VOCs were non-detectable except PCE which was detected at 0.95 µg/m³) and no chlorinated VOCs were detected in the March 17th sample. The ambient air data are summarized on **Table 3-1** and **Figure 12**.

3.6 Soil Investigation

The soil investigation during the RI was completed in several phases in conjunction with the monitoring well installation program to assess the extent of groundwater impacts. Each phase of soil investigation during the RI is summarized below.

3.6.1 Dry Cleaner Area Soil Investigation

Soil borings were advanced in the basement of the Dry Cleaner (SB-1, SB-2 and SB-3) in October 2015 to further characterize the nature and extent of contaminants in subsurface soils at the Site. Per the NYSDEC approved RIWP, soil borings SB-1 through SB-3 were advanced to 5.0 feet below the basement slab. Soil sample locations are detailed on **Figure 13** and analytical results are summarized in **Table 3-2** through **Table 3-5**. Analytical report summaries for all soil sampling events are provided in **Appendix E**.

Soil samples were collected from each boring location at the 1.0- and 5.0- foot bgs intervals (i.e., two samples per boring location, total of six samples) and submitted to SGS for analysis of Target Compound List (TCL) VOCs by EPA Method 8260C, TCL SVOCs by EPA Method 8270D, TCL pesticides by EPA Method 8081B, polychlorinated biphenyls (PCBs) by EPA Method 8082A, herbicides by EPA Method 8151 and Target Analyte List (TAL) metals by EPA Methods 6010C and 7471B.

For the purposes of this RIR, the soil analytical data are compared to the NYSDEC SCOs included in Part 375-5-6.8 - NYSDEC Remedial Program Soil Cleanup Objectives. As summarized in **Table 3-2**:

- PCE was detected in all six or the soil samples (both depths at all three boring locations) at concentrations ranging from 0.0183 to 0.0448 milligrams per kilogram (mg/kg), which are well below the NYSDEC Unrestricted Use SCO of 1.3 mg/kg for PCE.
- TCE was also detected in most of these soil samples at concentrations ranging from 0.004J to 0.0088J mg/kg. The NYSDEC Unrestricted Use SCO for TCE is 0.47 mg/kg.

• The only other TCL VOC detected was acetone at concentrations ranging from 0.003 to 0.0751 mg/kg (in the shallow soil sample collected from SB-2). The deeper sample analyzed from SB-2 contained acetone at a concentration of 0.0196 mg/kg. Acetone, which is a common laboratory artifact, has a NYSDEC Unrestricted Use SCO of 0.05 mg/kg and a Residential SCO of 100 mg/kg.

The results for SVOCs, pesticides, herbicides, PCBs, and inorganics (**Table 3-3** through **Table 3-5**) are summarized below:

- The only TCL SVOC detected in any of the six samples was bis(2-ethylhexyl)phthalate at concentrations ranging from 0.223 to 1.66 mg/kg. The NYSDEC has not assigned a SCO to this TCL SVOC.
- The only TCL pesticides detected in the samples were 4,4'-DDD at 0.0072 mg/kg and 4,4'-DDT at 0.0019 mg/kg. The NYSDEC Unrestricted Use SCOs for these TCL pesticides are 0.0033 mg/kg. No TCL PCBs or herbicides were detected in any of the six soil samples.
- The only TAL metals detected at concentrations exceeding their respective NYSDEC Unrestricted Use SCOs were iron at concentrations ranging from 9,740 mg/kg to 12,700 mg/kg (Supplemental Residential SCO of 2,000 mg/kg), cobalt in one sample at 32.2 mg/kg (NYSDEC Supplemental Residential SCO of 30mg/kg) and nickel ranging from 376.0 to 50.7 mg/kg (NYSDEC Unrestricted Use SCO of 30 mg/kg and Residential SCO of 140 mg/kg.
- Iron was detected in all soil samples collected from the basement subsurface at concentrations above the NYS residential SCO of 2,000 mg/kg which is not uncommon in urban areas of New York City. There are no commercial or groundwater protection standards for iron. Cobalt was detected in soil sample SB-2 at 1.0 feet below the basement slab at a concentration 32.3 mg/kg, just above the NYS residential SCO. Based on the Site history, it is unlikely that the cobalt and iron detected in the soil beneath the basement slab of the Dry Cleaner Area are related to historical Site uses. Additionally, based on the depth of these soil samples, a minimum of 9.0 feet bgs, the detections of cobalt and iron are most likely related to natural conditions in the subsurface.

3.6.2 Soil Investigation During Monitoring Well Installations

In January 2016, March 2016, and June 2016, soil samples were collected from the monitoring well boreholes during monitoring well installations to further define the nature and extent of groundwater impacts previously identified in the vicinity of the Dry Cleaner Area. Monitoring wells MW-11 and MW-4S were installed in January 2016. Monitoring wells MW-5, MW-6, and MW-7 were installed in March 2016. Monitoring wells MW-8, MW-9, and MW-10 were installed in June 2016. Monitoring well locations for all phases of investigation are indicated in **Figure 4**.

Soil samples were collected at the smear zone and below the water table during these installations to aid in evaluation of off-site impacts and contaminant mass distribution. The following soil samples were collected during well installation activities:

- Soil boring MW-11 was advanced to 50 feet bgs and soil samples we collected at 20 feet bgs, 45 feet bgs and 50 feet bgs.
- Soil boring MW-4S was advanced to a 40 feet bgs and soil samples were collected at 30 and 40 feet bgs.
- Soil boring MW-5 was advanced to 35 feet bgs and soil samples were collected at 22.5 feet and 35 feet bgs.
- Soil boring MW-6 was advanced to 30 feet bgs and soil samples were collected at 23 and 30 feet bgs.
- Soil boring MW-7 was advanced to 34 feet bgs and soil samples were collected at 23 and 34 feet bgs.

All soil samples during this phase of investigation were analyzed for VOCs via method SW846 8260C. Acetone was detected in the soil collected from soil boring MW-4S at 30 feet bgs at a concentration of 0.058 mg/kg and from soil

collected from soil boring MW-6 at 23 and 30 feet bgs at concentrations of .203 and .283 respectively. All acetone concentrations were detected above the NYS groundwater protection SCO; however, groundwater is a known laboratory contaminant and data should be interpreted cautiously. Groundwater in these two sample locations was encountered at 8.5 feet bgs in soil boring MW-4S and at approximately 20 feet bgs in soil boring MW-6 therefore, these concentrations of acetone detected in the soil above the NYS groundwater protection SCO are saturated and will continue to be monitored via groundwater sampling. No further soil delineation in relation to acetone is warranted.

To complete the delineation of PCE in groundwater, four additional monitoring wells were installed in June 2016. Soil samples were collected in these boreholes for further characterization of the subsurface soils in the area and the distribution of COC mass in area soils. Soil borings MW-8, MW-9 and MW-10 were advanced to 30 feet bgs and soil samples were collected at 19.5 and 30 feet bgs in soil borings MW-8 and MW-10 and at 19.4 and 30 feet bgs in soil boring MW-9. Soil boring MW-9. Soil boring MW-51 was advanced to 50 feet bgs and soil samples were collected at 20, 23, 31 and 50 feet bgs. The soil samples were submitted to a NYSDEC certified laboratory for analysis of TCL VOCs via method SW846 8260C. The analytical results reported that no COCs exceeded any of the NYS SCO from the soil borings MW-8, MW-9, MW-10 and MW-51.

Based upon the data at and immediately adjacent to the Dry Cleaner Area and the additional delineation further downgradient in conjunction with groundwater investigation, soil impacts have been delineated sufficiently to determine remediation needs, if any, for the Site.

3.7 Groundwater Investigation

The groundwater investigation at the Site was completed in phases with concurrence of NYSDEC so that data initially obtained could be used to optimize placement of future wells and sample locations and depths. This iterative process resulted in efficient definition of the chlorinated VOC plume both on-Site and at the adjacent off-site properties. The investigation approach and results are summarized in the following sections.

3.7.1 Initial Phase II Groundwater Investigation

In July 2012, three monitoring wells were installed onsite as part of a Phase II environmental site assessment (ESA). Monitoring wells MW-1 through MW-3 were installed to approximately 28 feet bgs. On July 27, 2012, the monitoring wells were surveyed to characterize groundwater flow. Groundwater was detected at an average of 20 feet bgs within the monitoring wells and groundwater elevation data collected from these three monitoring wells suggested groundwater flow was consistent with the regional flow direction which was to the south / southeast and south / southwest. Groundwater from the initial three monitoring wells was sampled on July 13, 2012. In addition to sampling the three permanent monitoring wells, groundwater samples were also collected from six temporary points B-1 through B-6 at this time. The groundwater samples were analyzed for VOCs and SVOCs. Analytical results were compared to the NYSDEC Class GA GWQS. PCE concentrations were detected in all nine groundwater samples above the applicable standards. TCE and *cis*-1,2 dichloroethene concentrations were detected in the groundwater samples, above their respective NYSDEC GA GWQS in the groundwater collected from monitoring wells, MW-1, MW-2 and temporary point B-1. No SVOCs were detected above the NYSDEC Class GA GWQS in any samples during this initial phase of investigation. See **Appendix A** for additional information regarding historic groundwater sampling events and results.

Based on the detections of chlorinated volatile organic carbons (CVOCs) above the NYSDEC Class GQ GWQS in the Site groundwater and the proximity to residential properties, further investigation to characterize and delineate the

COCs in groundwater was required. The additional investigations performed during the RI are summarized in the subsequent sections.

3.7.2 Delineation of Groundwater Impacts

The RIWP proposed to complete a MIP investigation to characterize and delineate the extent of the CVOC impacts in the surrounding soils. The MIP data was collected with the intent of confirming the final locations of the proposed monitoring wells. However, as discussed previously in **Section 3.4**, the MIP investigation did not indicate a high level of CVOCs in the areas assessed, even near monitoring well MW-1, a known location of CVOC impacts based upon initial sampling from "permanent" monitoring well located immediately downgradient of the suspected source area at the Dry-Cleaning Area. For this reason, a modified monitoring well installation plan was approved by the NYSDEC in a December 18, 2015, telephone conversation. Using a phased in approach, two intermediate groundwater monitoring wells were installed at the source near monitoring well MW-1 (new wells MW-11 and upgradient of the source near monitoring well MW-4². A full round of synoptic groundwater elevation data and COC analytical data was then used to determine the final delineation locations. The final monitoring well network is indicated in **Figure 4**. Monitoring well construction details for these three wells are summarized below and in **Appendix F**:

- Monitoring well MW-1I was installed on January 18 to 50 feet bgs. The monitoring well was constructed using two-inch diameter PVC with a 10-foot section of 0.010 slotted PVC screen set from 40 to 50 feet bgs.
- Monitoring well MW-4S and MW-4I were installed on January 19, 2016, in proximity to each other. Monitoring well MW-4S was installed west of the suspected source area to a total depth of approximately 27.5 bgs with 10-feet of two-inch diameter, 0.010 slotted PVC screen form 17.5- to 27.5-feet bgs. Monitoring well MW-4I was installed to 40 feet bgs with a10-foot, two-inch diameter section of 0.010 slotted PVC screen set from 30- to 40-feet bgs.

The groundwater was sampled on February 10, 2016, following the installation of these monitoring wells using the low flow purge method. The sampling sheets and field parameters collected during this monitoring and sampling event are included in **Appendix F**. The groundwater samples were analyzed for VOCs via USEPA method 624 and the groundwater samples collected from monitoring well MW-2 were also analyzed for SVOCs USEPA method 625, pesticides USEPA method 608, herbicides SW846 8151, and priority pollutant metals via PM 13. Sample results for groundwater sampling are summarized in **Tables 3-6** (VOCs), **Table 3-7** (SVOCs), **Table 3-8** (Pesticides and Herbicides), **Table 3-9** (Inorganics) and **Figure 14**. Laboratory analytical summary reports for groundwater samples are provided in **Appendix G**.

PCE, cis-1,2-dichloroethene, and TCE were detected in the groundwater samples above the NYSDEC Class GA GWQS in MW-1S (PCE at 740 µg/l, TCE at 9.7 µg/l, and DCE at 6.7 µg/l). PCE was also detected above the NYSDEC Class GA GWQS in the groundwater samples collected from monitoring well MW-2 (249 µg/l) and MW-3 (11.9 µg/l). No SVOCs, pesticides, herbicides, or metals were detected in monitoring well MW-2.

Based on the data from the February 10, 2016, groundwater monitoring and sampling event, Apex, as discussed with the NYSDEC, installed three additional monitoring wells to further delineate the COCs in groundwater. On March 9,

²Originally proposed MW-4 was renumbered MW-4S after installation of the intermediate well MW-4I to avoid confusion. Therefore, only wells MW-4S and MW-4I are indicated on figures for consistency. MW-4 and MW-4S are the same well.

2016, monitoring wells MW-5 through MW-7 were installed as indicated below:

- Monitoring wells MW-5 (later renamed MW-5S) and MW-6 were installed to 27 feet bgs with 10-feet of twoinch diameter, 0.010 slotted PVC screen from 17 to 27 feet bgs.
- Monitoring well MW-7 was installed to 30 feet bgs with 10-feet of two-inch, diameter, 0.010 slotted PVC screen from 20 to 30 feet bgs.

Groundwater was sampled from these monitoring wells on March 18, 2016 (see **Table 3-6** through **Table 3-9** and **Figure 14**). The groundwater samples were analyzed for VOCs via USEPA method 624 and the groundwater samples collected from monitoring well MW-5 were also analyzed for SVOCs USEPA method 625, pesticides USEPA method 608, herbicides SW846 8151, and priority pollutant metals via PM 13 per a NYSDEC request. Concentrations of cis-1,2-dichloroethene, trans-1,2-dichloroethene, PCE and TCE were detected above the NYSDEC Class GA GWQS in the groundwater sample collected from monitoring well MW-5/5S (PCE at 3,490 µg/l, TCE at 106 µg/l, and DCE at 687 µg/l). All VOCs were either below their applicable standard or non-detect in monitoring wells MW-6 and MW-7. No SVOCs, pesticides, herbicides, or metals were detected in monitoring well MW-5/5S.

Because COCs were detected in monitoring well MW-5, the most downgradient well during the March 2016 sampling event, four additional monitoring wells were installed in June 2016. In addition to horizontal plume delineation, Apex installed monitoring well MW-5I to characterize and delineate the COCs vertically at the MW-5 location where elevated COCs were detected in the shallow groundwater zone. On June 23, 2016 through June 24, 2017, the four monitoring wells (MW-5I, MW-9, MW-10, and MW-11) were installed as follows:

- Monitoring well MW-5I was installed to 50 feet bgs with 10-feet of two-inch diameter, 0.010 slotted PVC screen from 40 to 50 feet bgs.
- Monitoring well MW-9 was installed to 27 feet bgs with 10-feet of two-inch diameter, 0.010 slotted PVC screen from 17 to 27 feet bgs.
- Monitoring wells MW-8 and MW-10 were installed to 27 feet bgs with 10-feet of two-inch diameter, 0.010 slotted PVC screen from 17 to 27 feet bgs.

These wells completed the installation of groundwater monitoring points.

Groundwater was collected for sampling from these monitoring wells on July 8, 2016. The groundwater samples were submitted to a NYSDEC certified laboratory and analyzed for VOCs via USEPA method 624 (see **Table 3-6**, **Figure 14**, and **Appendix G**). PCE was detected in the groundwater from monitoring wells MW-5I at 18.7 µg/L, MW-9 at 7.5 µg/L and MW-10 at 17.6 µg/L. No other COCs were detected above the NYSDEC Class GA GWQS.

Based on the results from the phased approach to groundwater characterizations and delineation, the CVOCs in the groundwater are delineated. As can be seen on **Figure 14**, there is a decreasing concentration gradient as groundwater moves away from the Site. Concentrations to the east and south of the Site are below the NYSDEC Class GA GWQS and monitoring well MW-8 to the southwest had no detections of CVOCs. Additionally, TCE and cis-1,2 DEC are degradation byproducts of PCE and these COCs were detected in groundwater samples collected at this Site indicating that natural biodegradation is taking place.

3.8 Summary of Data Usability

This section provides a summary of RI data quality based on usability toward meeting data quality objectives (DQOs) for the Site. The full Data Usability Summary Reports (DUSRs) is provided in **Appendix H** (under separate cover). The Data Usability Report was completed by Kelsi Evans, a chemist out of Apex's Portland., Oregon office. Ms. Evans'

resume and qualification are included in **Appendix H**. Ms. Evans was not involved in any aspect of the RI project other than the independent data validation.

Based on the data validation, the data met DQOs and are of sufficient quality to meet the objectives of the RI.

3.9 Investigation-Derived Waste Disposal

Investigation derived waste (IDW) was managed as proposed in the NYSDEC approved RIWP. IDW was containerized in 55-gallon steel drums and staged onsite for appropriate characterization and disposal following the procedures for IDW management outlined in the Field Sampling Plan. Personal Protective Equipment (PPE) and spent disposable sampling materials were segregated and placed in NYS Department of Transportation (DOT)-approved 55-gallon steel drums. Decontamination water and monitoring well purge water was stored in DOT approved 55-gallon steel drums. Waste storage containers were appropriately labeled with the contents, generator, location, and date and properly secured at the site for subsequent off-site transportation and disposal by the contractor.

One representative sample was collected from the solid IDW (i.e., drill cuttings) and one representative sample was collected from the liquid IDW (i.e., development water/monitoring well purge water) generated by the field activities. The samples were submitted to the laboratory for analysis of the parameters required by the off-site disposal facility. Apex used the analytical results from the waste characterization samples to profile the IDW for proper disposal.

4.0 Nature and Extent of COCs and Conceptual Site Model

This section discusses the analytical data for samples of soil, groundwater, and exterior soil vapor collected at the Site from the perspective of delineation and potential for migration.

4.1 Nature and Extent of COCs in Soil

This section provides an evaluation of the nature and extent of COCs in soil. The data obtained from soil sampling met the RI goal of determining the extent of COCs and was sufficient for determining the nature and extent of COCs in soil at the Site.

Based upon the discussion of **Section 3.6**, the analytical results indicate that a few VOCs were detected in soil above SCGs. The primary VOCs that were detected in soil above SCGs include acetone, PCE, TCE and cis-1,2 dichloroethene. These COCs exceeded the NYSDEC SCO for protection of groundwater (as is discussed in **Section 4.2**). Acetone was detected in soil sample SB-2 at 1.0 feet bgs, MW-4S at 30 feet bgs, and MW-6 at 23 and 30 feet bgs. However, acetone is a common laboratory artifact and is not considered a COC with respect to an on-site source. PCE and cis-1,2 DCE were detected above the groundwater protection SCO in soil samples MW-5 at 30 and 35 feet bgs which is below the water table and likely related to groundwater impacts in that area. TCE was detected above the groundwater protected at that location and depth, is likely present due to elevated groundwater concentrations in that area.

The soil data collected during the RI indicated that soil impacts have been delineated and additional soil investigation activities are not warranted.

4.2 Nature and Extent of COCs in Groundwater

This section provides an evaluation of the nature and extent of COCs in groundwater. The data obtained from sampling groundwater collected from the monitoring wells that were installed for the RI met the RI goal of determining the nature and extent of COCs in groundwater. An analysis of the COCs in groundwater is presented below.

The analytical results indicate that several VOCs (PCE and its degradation products TCE and cis-1,2 DCE primarily) were detected in groundwater above their respective SCGs. The following is a summary of the chlorinated VOC distribution in groundwater based upon the most recent groundwater sampling events in 2016:

- PCE was detected in all shallow groundwater monitoring wells except for MW-8. The highest PCE concentration in the shallow zone was identified off-site at MW- 5S (3,490 µg/l PCE) which is located approximately 100 feet downgradient of the on-site dry-cleaner area. Other than wells MW-1S (740 µg/l PCE), MW-2S (249 µg/l PCE), and MW-5S (3,490 µg/l PCE) which are located immediately near and downgradient of the dry-cleaner area, PCE concentrations in all remaining shallow groundwater monitoring wells in the study area ranged from non-detectable at MW-8 to 17.6 µg/l at MW-10 (approximately 300 feet downgradient of the dry-cleaner area). These data support that elevated PCE concentrations in the shallow zone are relatively local to the apparent source area. This conclusion is further supported by the fact that upgradient PCE concentrations in the shallow zone at MW-3 were 11.9 µg/l which is similar to most PCE concentrations outside of the apparent source area and immediately downgradient.
- PCE concentrations as the depth increased from grade were also evaluated during the RI based upon MIP data screening. The highest PCE concentration in the intermediate zone was 18.7 at MW-51 (located in the

same location as MW-5S). There was a two order of magnitude decline in PCE concentrations from the shallow zone to the intermediate zone at MW-5S/MW-5I.

TCE and 1DCE concentrations were also the highest at MW-5S (106 μg/l TCE and 687 μg/l DCE). The concentrations of TCE and DCE in all other wells (both shallow and intermediate) was under 10 μg/l for both parameters with most locations under 5 μg/l.

As discussed in **Section 3.5** and indicated in **Figure 14**, the distribution of CVOC concentrations in groundwater has been delineated horizontally and vertically. There appears to be a relatively narrow band of COC impacts in shallow groundwater originating at the Dry Cleaner Area (near MW-1S) and extending along the western side of Bay 25th Street to the MW-5S area before declining rapidly in concentration toward MW-9 and MW-10. Only minimal impacts were noted both east and west of this narrow band of shallow groundwater impacts. Intermediate depth groundwater impacts at MW-11 were non-detectable indicating minimal to no significant vertical migration from MW-1S impacts. Similarly, the PCE concentration at MW-51 is several orders of magnitude lower than the MW-5S concentrations indicating that vertical migration is defined.

The data obtained during the RI have defined the nature and extent of groundwater warranting remediation and / or further monitoring.

4.3 Nature and Extent of COCs in Soil Vapor and Ambient Air

This section provides an evaluation of the nature and extent of COCs in sub slab soil vapor and indoor air quality at the Site (Block 6467, Lot 1), at the adjacent property, 8831 and 8841 20th Avenue (Block 6467, Lot 12) and the residential property to the southeast of the Site, 2036 Cropsey Avenue (Block 6469, Lot 1). The data obtained from soil vapor sampling met the RI goal of determining the nature and extent of CVOC-related COCs in soil vapor at the Site. An analysis of COCs in soil vapor is presented below.

On-Site Vapor and Indoor Air:

- The highest concentrations of chlorinated VOCs in soil vapor were detected on-site during the July 2012 soil vapor screening event with 420,000 µg/m³ of PCE and 6,600 µg/m³ of TCE detected in SV-1 underlying the dry-cleaning establishment. On-site soil vapor concentrations west of the dry-cleaning area and further to the west off-site were less than 76 µg/m³ PCE and less than 1.5 µg/m³ TCE. Indicating localized, maximum concentrations (see Figure 12 and Table 3-1).
- The highest indoor air concentrations of PCE were detected at SV-2 (in the dry-cleaning establishment) in the July 2012 sampling event (100 μ g/^{m3} PCE). The highest indoor air concentrations for TCE were detected in this same location in the July 2012 sampling event (18 μ g/^{m3} TCE).

Off-Site Vapor and Indoor Air:

- The highest off-site soil vapor impacts were noted in SV-3 through SV-5 underlying the off-site parking garage located approximately 50 feet to the south-southwest (downgradient) of the apparent dry-cleaner source area. The maximum off-site concentrations of PCE ranged from 140,000 to 210,000 µg/m³ in SV-3 through SV-5 in July 2012. Maximum TCE concentrations in this area during the July 2012 sampling event ranged from 580 to 790 µg/m³.
- In the most recent soil vapor and indoor air quality sampling event in April 2017, sub-slab vapor concentrations of PCE at SV-3, SV-4, and SV-5 were 82,700 µg/m³, 8,070 µg/m³, and 6,520 µg/m³, respectively). These concentrations were one to two orders of magnitude lower than the 2012 sampling results at these same locations. TCE sub-slab concentrations at SV-3 and SV-4 were the same order of

magnitude as the past 2012 sampling event and the concentrations of TCE in sub-slab vapor at SV-5 dropped to non-detectable levels.

- Indoor air sample concentrations at SV-3, SV-4, and SV-5 for PCE were on the same order of magnitude in the 2017 sampling event as in the past 2012 sampling event with slight increases noted at all three locations (minimum of 1.9 µg/m³ and maximum of 9.5 µg/m³ for PCE over both events). TCE concentrations in indoor air samples at SV-3, SV-4 and SV-5 were all an order of magnitude lower in the 2017 sampling event as compared to the 2012 sampling event with a maximum concentration of 1.1 µg/m³TCE detected during the July 2012 event and a maximum of 0.39 µg/m³TCE detected during the 2017 sampling event.
- Off-site soil vapor and indoor air concentrations for COCs, including the upgradient locations north and northeast of the site, decreased sharply with distance from the maximum concentration locations identified above indicating a relatively focused area of maximum concentrations that is generally consistent with the groundwater plume maximum concentration area. Although the concentrations of upgradient locations north and northeast of the site declined rapidly, PCE and TCE soil vapor detections are present upgradient of the site.

To address the concentrations of chlorinated VOCs discussed above for the dry-cleaner / shopping center area and the adjacent residential parking garage areas, two mitigation systems were employed. An IRM was proposed and installed for a SSDS at the dry cleaner facility / retail shopping center where the highest concentrations of PCE and TCE were detected in both sub-slab concentrations and in indoor air samples. In addition, below grade parking garage fan upgrades were also been installed at the adjacent residential building parking garages to prevent vapors from entering the occupied spaces of the facility at levels above the parking garages.

Both the parking garage fan systems and the SSDS underlying the shopping center are designed to run 24 hours a day / 7 day a week. The parking garage system details are provided in **Appendix I**. The shopping center SSDS system details are provided separately in the Interim Remedial Measures (IRM) Installation and Start-up Report (Apex, January 2018). Both systems operate to create a pressure differential between occupied areas and unoccupied areas / the subsurface. For the parking garage fan systems, a draft / vacuum is induced in the sub-grade parking garage so that the pressure in overlying, occupied structures is higher than in the garage area. This pressure differential will prevent contaminant vapors underlying the building from entering occupied spaces above the subsurface parking garage. The SSDS at the shopping center works on the same principle with the below grade vapor extraction points creating a subsurface vacuum relative to the overlying, occupied areas of the shopping center.

Additional monitoring for both mitigation systems will be ongoing and coordinated with NYSDEC.

4.4 Conceptual Site Model

This section develops and discusses the CSM for the Site. Previous sections summarized relevant Site conditions, including the Site history and usage, geology and hydrogeology, and distribution of COCs in the environment (i.e., quality of soil, groundwater, soil vapor and indoor). This CSM relates current conditions to the historical dry cleaner operations at the Site. The purpose of the CSM is to more specifically explain the source of the COCs and the movement of the COCs through the subsurface. The data collected during the RI was used to draw conclusions as to the nature and extent of the COCs. The goal of the CSM is to understand, to the extent it can be, how and why COCs entered the soil, groundwater, and soil vapor and where COCs will migrate if not remediated. The following is a summary of the CSM based upon the available data. The CSM will be continually re-evaluated as additional data are available to ensure protection of human health and the environment.

• PCE was found to be the primary COC in soil vapor, soil and groundwater at the Site. The likely source of the PCE is the operations of a former dry-cleaning tenant at the location currently occupied by GLY Cleaners, an

unrelated dry-cleaning operation. Dry cleaner operations reportedly have not used PCE since 2003; however, based on file reviews, PCE was used at the facility prior to 2003. The current dry cleaner tenant began operation in 2007 and has not reportedly used PCE. It also should be noted that PCE (and its degradation products) has been detected upgradient as well so there is likely some comingling of PCE from an independent upgradient source.

- The specific release(s) that occurred at this Site are unknown. Based on the historic Site use as a dry cleaner and the concentrations of CVOCs in the subsurface under the basement slab and directly downgradient of the Site, it is likely that dry cleaning solvents historically spilled and entered the subsurface either directly through the basement floor (i.e., via cracks or floor deterioration or historical drains that may have been present). Given impacts in the rear "yard" behind the dry cleaner space, surface releases also may have been possible. These localized impacts are the likely source of soil vapors that are present at the Site.
- As discussed in **Section 4.2**, COCs entering the soil likely migrated primarily vertically into the groundwater underlying the Dry Cleaner Area and then were transported primarily horizontally with shallow groundwater flow to the south-southeast (off-site). Vertical migration appears to have been minimal once COCs entered the water table based upon the available data. There may be a southwesterly flow component for the relatively low concentrations of COCs detected in the intermediate groundwater zone.
- Groundwater migration appears to be occurring in a very narrow band from the Site along the western side of Bay 25th Street toward MW-5/5S and ultimately toward MW-9 and MW-10. There is a decrease in concentration of COCs by several orders of magnitude at MW-9 and MW-10 in comparison to the source area wells at MW-1 and the MW-5 wells indicating that the plume area is localized.
- Soil vapor data is generally consistent with the groundwater plume (i.e., highest concentrations underlying the apparent source area at the dry-cleaning location and immediately downgradient) that has been defined and will likely require management until groundwater impacts are remedied. The IRMs implemented to date and to be implemented can effectively manage soil vapor concerns.

5.0 Soil Vapor Mitigation Measures and IRMs

Per the NYSDEC-approved IRM work plan, the following work components were conducted as part of soil vapor mitigation IRMs.

5.1 Shopping Center SSDS IRM

A Sub-Slab Depressurization System (SSDS) pilot test was completed in March 2016 to confirm the applicability of SSDS to mitigate sub-slab soil vapors migration into occupied spaces above grade. On March 30 and 31, 2016, three vacuum points and five vacuum monitoring points were installed in preparation for the SSDS tests (see **Figure 15**). Three tests were completed to assess the radius of influence (ROI) of vacuum applied to the vacuum points. The results of the pilot tests indicated that the lateral extent of the vacuum field was extended to underlie the entire Site, including each existing tenant space and between tenant spaces with a maximum ROI of 78 feet from the vacuum source (**Figure 16**). The NYSDEC accepted pilot test data and approved the IRM Work Plan for system installation based on successful pilot testing and system installation is scheduled for September 2017. Additional information is available in the NYSDEC-approved IRM Work Plan under separate cover.

Following successful pilot testing the SSDS IRM was installed. The system is currently operational. Additional information regarding the SSDS IRM including construction as-built drawings and start-up data are provided in the SSDS IRM Installation Completion and Start-up Report (Apex - March, 19, 2018).

5.2 Residential Building Garage Ventilation System Upgrades

Due to the presence of elevated concentrations of the COCs in soil vapors underlying the site and the adjacent residential buildings, a mitigation measure consisting of parking garage ventilation system upgrades was implemented. The garage ventilation system IRM details are included in **Appendix I**.

6.0 Human Health Exposure Assessment

In accordance with the NYSDEC DER-10 Technical Guidance for Site Investigation and Remediation (NYSDEC, 2010), a qualitative exposure assessment was performed and a Human Health Exposure Assessment (HHEA) was completed. The HHEA identified COCs, identified potentially complete exposure pathways, and evaluated the potential for exposure of human receptors to Site-related COCs assuming no further remedial actions were conducted.

Soil, groundwater, and ambient air associated with the Site do not present potentially complete exposure pathways for commercial workers, consumers, or residents based on current land use and are not anticipated to represent complete future exposure pathways for these receptors unless there is construction or maintenance activities that would break the existing slab at the Site or the surface coverings that currently overlie the groundwater COC plume. Future construction worker scenarios will be addressed in a future Site Management Plan (SMP).

Soil vapor beneath the Site and the adjacent property to the south presents a potential for a completed pathway to indoor air and as stated within this report, health and safety measures; a sub-slab depressurization system and ventilation system upgrades are being implemented to mitigate the potential risk.

| Environmental Media & Exposure Route | Human Exposure Assessment (examples) |
|---|---|
| Direct contact with surface soils (and incidental ingestion) | There are no surface soil impacts at this Site. There is only one limited area of soil impacts underlying the basement slab of the Dry-Cleaning Area. Under current operating conditions, there is no completed exposure pathway. If current site conditions change or if construction activities require demolition of the basement floor slab, additional measures may be required to ensure worker protection. Contingent protection measures for future site use scenarios can be addressed in a Site Management Plan to be developed. |
| Direct contact with subsurface soils (and incidental ingestion) | The soil impacts at this Site are limited to exceedances of the groundwater protection standard not the direct contact standard. However, there is one soil sample location impacted above the respective health based NYSDEC SCO. People are not coming into contact because soil impact is below the basement slab of the drycleaner. People are not coming into contact with the other Site soil impacts because they are > 23 feet bgs and are under pavement or landscaping Contingent protection measures for future site use scenarios can be addressed in a Site Management Plan to be developed. |
| Ingestion of groundwater | Contaminated groundwater is not being used for drinking water, as the area is served by the public watersupply. There are no known domestic water supply wells in thearea. People can ingest GW if private wells are installed on the property or under possible dewatering scenarios. These potential future impacts would be addressed in an SMP to be developed. |

Per the NYSDEC requirements, the following summary of current and potential exposure table is included.

| Direct contact with groundwater | People can come into contact with groundwater if they complete ground- intrusive work at the Site or the adjacent underground parking garage facilities. However, the average depth to groundwater is 20 feet bgs so it is unlikely at shallow depths. Contingent protection measures for future site use scenarios can |
|---|--|
| | be addressed in a Site Management Plan to be developed. |
| Inhalation of air (exposures related to soil vapor intrusion) | A ventilation system fan upgrade has been installed in the parking garages of the offsite buildings to prevent the indoor air quality from being affected by the sub-slab vapor impacts. |
| | • A sub-slab depressurization system is being installed in the on-site shopping center to prevent the indoor air quality from being affected by the sub-slab impacts. |
| | Post-installation indoor air samples will be collected from the on-site building to verify that additional actions are not needed to address exposures related to soil vaporintrusion. |
| | • The SMP to be developed will address these issues further. |
| | |

7.0 Findings and Conclusions

The soil, soil vapor, and groundwater data evaluated during this RI have resulted in delineation of the nature and extent of COC impacts. There is sufficient data to evaluate potentially applicable remedial measures to address the impacts to media described within this RI. The following is a summary of the RI conclusions:

- The primary COCs in soil vapor are chlorinated organic compounds, notably PCE and its degradation compound TCE. The likely, but unconfirmed, source of soil vapor impacts are the historic dry cleaner operations within the 2002-20024 shopping center. As indicated in the CSM, releases of dry cleaning solvents likely entered the subsurface at the Dry-Cleaning Area and migrated vertically into groundwater underlying the Site. Once COCs were in groundwater, they migrated to the south-southeast (off-site) with groundwater flow resulting in soil vapor impacts in overlying areas. Migration of soil vapor impacts are generally consistent with groundwater plume impacts. COCs are present underlying the shopping center as well as the adjacent, off-site parking garages. The most significant soil vapor impacts emanate from the Site along the west side of Bay 25th Street and extending off-site to the vicinity of MW-5/5S. COC concentrations decrease several orders of magnitude beyond MW-5/5S and also to the east and west beyond approximately 100 feet laterally from the areas exhibiting the most significant groundwater impacts
- The primary VOCs that were detected in soil above SCGs include acetone, PCE, TCE and cis-1,2 dichloroethene. These COCs exceeded the NYSDEC SCO for protection of groundwater. Unsaturated zone impacts are very localized immediately underlying the Dry Cleaner Area. Acetone was detected in soil sample SB-2 at 1.0 feet bgs, MW-4S at 30 feet bgs, and MW-6 at 23 and 30 feet bgs. However, acetone is a common laboratory artifact and is not considered a COC with respect to an on-site source. PCE and cis-1,2 DCE were detected above the groundwater protection SCOs in soil samples MW-5 (off-site) at 30 and 35 feet bgs which is below the water table and likely related to groundwater impacts in that area. TCE was detected above the groundwater protection SCO in soil sampled during MW-5 installation at 35 feet bgs and similar to PCE and cis-1,2 DCE detected at that location and depth, is likely present due to elevated groundwater concentrations in that area. Soil impacts are delineated.
- Groundwater impacts are present. Analytical results indicate that several VOCs (PCE and its degradation products TCE and cis-1,2 DCE primarily) were detected in on-site and off-site groundwater above their respective SCGs. The distribution of CVOC concentrations in groundwater has been delineated horizontally and vertically. There appears to be a relatively narrow band of COC impacts in shallow groundwater originating at the Dry Cleaner Area (near MW-1S) and extending along the western side of Bay 25th Street to the MW-5S area before declining rapidly in concentration toward MW-9 and MW-10. Only minimal impacts were noted both east and west of this narrow band of shallow groundwater impacts. Intermediate depth groundwater impacts at MW-11 were non-detectable indicating minimal to no significant vertical migration from MW-1S impacts. Similarly, the PCE concentration at MW-51 is several orders of magnitude lower than the MW-5S concentrations indicating that vertical migration is defined.

In summary, the RI for this Site is complete and of sufficient scope and detail to allow evaluation of potential remedial alternatives to address environmental impacts. The following general path forward is recommended for the site:

- Develop a Remedial Action Plan (RAP) for the Site which will include a Feasibility Study / Alternatives Evaluation to screen potentially applicable remedial technologies and select the most appropriate remedial approach for the Site.
- As part of the RAP development, a Pre-Design Investigation (PDI) will be implemented. The main intent of the PDI will be to update the baseline conditions of the site in advance of remedial design and to obtain data necessary to support final remedy selection. Due to the phased nature of the RI, it is important to complete a comprehensive groundwater sampling event(s) which includes all final wells installed to serve as the basis for engineering evaluations.

- Depending upon the remedies considered for implementation, bench-scale treatability studies and / or pilot tests may be required before final design. These studies would be coordinated with the RAP development and detailed Remedial Design (RD) schedule with NYSDEC.
- Following the detailed remedial design, the selected remedial approach would be implemented and the appropriate engineering and institutional controls, Site Management Plan (SMP), etc. would be developed and completed.

Tables

<u>Table 3-1</u> Soil Vapor and Indoor Air Sampling Results Summary

| | Historical NYSDOH Regulations | | 2013 and 2015 NYSDOH Revised Standards | I Initial screening sampling | | | March 2016 sampling event including sub-slab and IAQ samples | | | | | | | | | | | | | April 2017 sampling event including sub-slab and IAQ samples | | | | | | |
|--|----------------------------------|----------------------------|---|------------------------------|------------------|------------------|--|------------------|------------------|--------------------|-----------------|--------------------|--------------------|------------------|------------------|------------------|--------------------|--------------------|-----------------|--|-----------------|--------------------|--------------------|--|--|--|
| Analyte Method TO-15 (ug/m ³) | Matrix 1 & 2 Sub-Slab Vapor | Matrix 1 & 2 Indoor Air | Air | SV-6 10/30/15 | SV-7 10/30/15 | SV-8 10/30/15 | SV-9* 3/4/16 | IAQ-9* 3/4/16 | SV-10 3/9/16 | IAQ-10 3/9/16 | SV-11 3/4/16 | IAQ-11 3/4/16 | Ambient 3/9/16 | SV-12 3/17/16 | SV-13 3/17/16 | SV-14 3/17/16 | Ambient 3/17/16 | IAQ-3 4/14/17 | SV-4 4/14/17 | IAQ-4 4/14/17 | SV-5 4/14/17 | IAQ-5 4/14/17 | Ambient 4/14/17 | | | |
| Acetone | NA | NA | | 634 | 1220 | 632 | 363 | 41.1 | 527 | 16 | 2.2 U | 9.7 | 13 | 10 U | 447 | 1310 | 2 | 0.086 U | 52 | 15 | 822 | 21 | 5.2 | | | |
| 1,3-Butadiene | NA | NA | | 0.27 U | 0.27 U | 0.27 U | 0.062 U | 1.7 J | 0.24 U | 0.062 U | 1.6 U | 0.55 | 0.062 U | 7.3 U | 0.24 U | 0.24 U | 0.062 U | 5.5 | 2.4 U | 0.062 U | 2.4 U | 5.8 | 0.062 U | | | |
| Benzene | NA | NA | | 4.8 | 5.4 | 6.1 | 0.64 | 6.4 | 3.5 | 2.6 | 89.1 | 8 | 1.1 | 33.9 J | 15 | 16 | 0.099 U | 30 | 4.2 U | 26 | 4.2 U | 30 | 0.93 | | | |
| Bromodichloromethane | NA | NA | | 0.87 U | 0.87 U | 0.87 U | 0.26 U | 1 U | 1 U | 0.26 U | 6.6 U | 0.26 U | 0.26 U | 31 U | 1 U | 1 U | 0.26 U | 0.26 U | 10 U | 0.26 U | 10 U | 0.26 U | 0.26 U | | | |
| Bromoform | NA | NA | | 0.85 U | 0.85 U | 0.85 U | 0.17 U | 0.65 U | 0.65 U | 0.17 U | 4.1 U | 0.17 U | 0.17 U | 20 U | 0.65 U | 0.65 U | 0.17 U | 0.17 U | 6.5 U | 0.17 U | 6.5 U | 0.17 U | 0.17 U | | | |
| Bromomethane | NA | NA | | 0.34 U | 0.34 U | 0.34 U | 0.07 U | 0.29 U | 0.29 U | 0.07 U | 1.8 U | 0.07 U | 0.07 U | 8.5 U | 0.29 U | 0.29 U | 0.07 U | 0.07 U | 2.9 U | 0.07 U | 2.9 U | 0.07 U | 0.07 U | | | |
| Bromoethene | NA | NA | | 0.35 U | 0.35 U | 0.35 U | 0.079 U | 0.32 U | 0.32 U | 0.079 U | 2.1 U | 0.079 U | 0.079 U | 9.6 U | 0.32 U | 0.32 U | 0.079 U | 0.079 U | 3.2 U | 0.079 U | 3.2 U | 0.079 U | 0.079 U | | | |
| Benzyl Chloride | NA | NA | | 0.52 U | 0.52 U | 0.52 U | 0.14 U | 0.57 U | 0.57 U | 0.14 U | 3.6 U | 0.14 U | 0.14 U | 16 U | 0.57 U | 0.57 U | 0.14 U | 0.14 U | 5.7 U | 0.14 U | 5.7 U | 0.14 U | 0.14 U | | | |
| Carbon disulfide | NA | NA | | 0.34 U | 0.34 U | 0.34 U | 2.2 | 0.4 U | 0.4 U | 0.097 U | 7.8 J | 0.097 U | 0.097 U | 12 U | 6.2 | 5.9 | 0.097 U | 0.31 J | 4 U | 0.097 U | 38.9 | 0.097 U | 0.097 U | | | |
| Chlorobenzene | NA | NA | | 0.6 U | 0.6 U | 0.6 U | 0.26 U | 1 U | 1 U | | 0.1 0 | 0.26 U | 0.26 U | 30 U | 1 U | 1 U | 0.26 U | 0.26 U | 10 U | 0.26 U | 10 U | | 0.26 U | | | |
| Chloroethane | NA | NA | | 0.23 U | 0.23 U | 0.23 U | 0.095 U | 0.37 U | 0.37 U | | 14 | 0.095 U | 0.095 U | 11 U | 0.37 U | 0.37 U | 0.095 U | 0.095 U | 3.7 U | 0.095 U | 3.7 U | 0.095 U | 0.095 U | | | |
| Chloroform | NA | NA | | 40 | 0.59 U | 12 | 1.1 | 0.32 U | 0.32 U | 0.083 U | 2.1 U | 0.083 U | 0.083 U | 9.8 U | 57.6 | 17 | 0.083 U | 0.083 U | 3.2 U | 0.54 J | 160 | 0.73 J | 0.083 U | | | |
| Chloromethane | NA | NA | | 0.25 U | 0.25 U | 0.25 U | 0.11 U | 1.4 J | 0.43 U | | 2.7 U | | 1.8 | 13 U | 0.43 U | 0.43 U | 0.39 J | 1.2 | 4.3 U | 1.4 | 4.3 U | | 1.2 | | | |
| 3-Chloropropene | NA | NA | | 0.34 U | 0.34 U | 0.34 U | 0.085 U | 0.34 U | 0.34 U | | 2.1 U | 0.085 U | 0.085 U | 9.7 U | 0.34 U | 0.34 U | 0.085 U | 0.085 U | 3.4 U | 0.085 U | 3.4 U | 0.085 U | 0.085 U | | | |
| 2-Chlorotoluene | NA | NA | | 0.67 U | 0.67 U | 0.67 U | 0.088 U | 0.35 U | 0.35 U | | 2.2 U | 0.088 U | 0.088 U | 10 U | 0.35 U | 0.35 U | 0.088 U | 0.088 U | 3.5 U | 0.088 U | 3.5 U | 0.088 U | 0.088 U | | | |
| Carbon tetrachloride | >250 Mitigate | >5 Mitigate | | 0.62 U | 0.62 U | 0.62 U | 0.2 U | 0.75 U | 0.75 U | | ° ° | 0.2 U | 0.6 J | 23 U | 0.75 U | 0.75 U | 0.2 U | 0.2 U | 7.5 U | 0.2 U | 7.5 U | | 0.2 U | | | |
| Cyclohexane | NA | NA | | 2.1 J | 2.6 J | 2.4 J | 0.96 | 0.22 U | 0.22 U | | 2.1.0 | 1 | 0.055 U | 396 | 10 | 7.9 | 0.055 U | 4.1 | 2.2 U | 3.8 | 2.2 U | | 0.055 U | | | |
| 1,1-Dichloroethane | NA | NA | | 0.49 U | 0.49 U | 0.49 U | 0.061 U | 0.25 U | 0.25 U | | 1.6 U | 0.061 U | 0.061 U | 7.3 U | 0.25 U | 0.25 U | 0.061 U | 0.061 U | 2.5 U | 0.061 U | 2.5 U | 0.061 U | 0.061 U | | | |
| 1,1-Dichloroethylene | NA | NA | | 0.44 U | 0.44 U | 0.44 U | 0.083 U | 0.33 U | 0.33 U | | 2.1 U | 0.083 U | 0.083 U | 9.9 U | 0.33 U | 0.33 U | 0.083 U | 0.083 U | 3.3 U | 0.083 U | 3.3 U | 0.083 U | 0.083 U | | | |
| 1,2-Dibromoethane | NA | NA | | 1.1 U | 1.1 U | 1.1 U | 0.32 U | 1.3 U | 1.3 U | 0.32 U | 8.5 U | 0.32 U | 0.32 U | 38 U | 1.3 U | 1.3 U | 0.32 U | 0.32 U | 13 U | 0.32 U | 13 U | 0.32 U | 0.32 U | | | |
| 1,2-Dichloroethane | NA | NA | | 0.4 U | 0.4 U | 0.4 U | 0.073 U | 0.28 U | 0.28 U | | 1.8 U | 0.073 U | 0.073 U | 8.5 U | 0.28 U | 0.28 U | 0.073 U | 0.073 U | 2.8 U | 0.073 U | 2.8 U | 0.073 U | 0.073 U | | | |
| 1,2-Dichloropropane | NA | NA | | 0.92 U | 0.92 U | 0.92 U | 0.1 U | 0.41 U | 0.41 U | 0.1 U | 2.6 U | 0.1 U | 0.1 U | 12 U | 0.41 U | 0.41 U | 0.1 U | 0.1 U | 4.1 U | 0.1 U | 4.1 U | 0.1 U | 0.1 U | | | |
| 1,4-Dioxane | NA | NA | | 0.9 U | 0.9 U | 0.9 U | 0.16 U | 0.65 U | 0.65 U | 0.16 U | 4.3 U | 0.16 U | 0.16 U | 19 U | 0.65 U | 0.65 U | 0.16 U | 0.16 U | 6.5 U | 0.16 U | 6.5 U | 5.4 | 0.16 U | | | |
| Dichlorodifluoromethane | NA | NA | | 2.8 J | 2.6 J | 3 J | 2.7 | 2.4 J | 2.8 J | 3.2 | 2.4 U | 2.8 | 3.4 | 11 U | 2.9 J | 2.9 J | 0.69 J | 2 | 3.8 U | 2 | 3.8 U | 1.9 | 2 | | | |
| Dibromochloromethane | NA | NA | | 1.4 U | 1.4 U | 1.4 U | 0.45 U | 1.8 U | 1.8 U | 0.45 U | 11 U | 0.45 U | 0.45 U | 53 U | 1.8 U | 1.8 U | 0.45 U | 0.45 U | 18 U | 0.45 U | 18 U | 0.45 U | 0.45 U | | | |
| trans-1,2-Dichloroethylene | NA NA | NA 100 Mitianto | | 0.32 U | 0.32 U | 6.3 | 0.11 U | 0.44 U | 0.44 U | | 18 J | 0.11 U | 0.11 U | 13 U | 0.44 U | 0.44 U | 0.11 U | 0.39 J | 4.4 U | 0.4 J | 4.4 U | 0.38 J | 0.38 J | | | |
| cis-1,2-Dichloroethylene | >1000 Mitigate | >100 Mitigate | | 0.39 U 0.64 U | 0.39 U 0.64 U | 6.3 0.64 U | 0.083 U 0.068 U | 0.33 U 0.28 U | 0.33 U 0.28 U | 0.083 U 0.068 U | 2.1 U | 0.083 U 0.068 U | 0.083 U 0.068 U | 120 8.2 U | 0.33 U 0.28 U | 0.33 U 0.28 U | 0.083 U 0.068 U | 0.083 U 0.068 U | 128 2.8 U | 0.083 U 0.068 U | 3.3 U 2.8 U | 0.083 U 0.068 U | 0.083 U 0.068 U | | | |
| cis-1,3-Dichloropropene | NA NA | NA NA | | 0.64 U | 0.64 U | 0.64 U | 0.068 U 0.12 U | 0.28 U | 0.28 U | 0.068 U | 1.8 U 3 U | 0.068 U | 0.068 U 0.12 U | 8.2 U | 0.28 U 0.47 U | 0.28 U | 0.000 U | 0.068 U | 2.8 U 4.7 U | 0.068 U | 2.8 U 4.7 U | 1.3 | 0.088 U 0.12 U | | | |
| m-Dichlorobenzene o-Dichlorobenzene | NA | NA | | 0.66 U | 0.66 U 0.72 U | 0.66 U | 0.12 U 0.096 U | 0.47 U | 0.47 U | 0.12 U | | 0.12 U 0.096 U | 0.12 U 0.096 U | 14 U | 0.47 U | 0.47 U | 0.096 U | 0.12 U 0.096 U | 4.7 U 3.8 U | 0.12 U 0.096 U | 4.7 U | 1.5 | 0.12 U 0.096 U | | | |
| p-Dichlorobenzene | NA | NA | | 0.46 U | 0.46 U | 0.46 U | 0.090 U | 0.66 U | 0.66 U | | | 0.090 U | 0.090 U | 19 U | 0.66 U | 0.66 U | 0.16 U | 0.090 U | 6.6 U | 0.090 U | 6.6 U | 2 | 0.16 U | | | |
| trans-1.3-Dichloropropene | NA | NA | | 0.40 U | 0.46 U | 0.40 U | 0.082 U | 0.33 U | 0.33 U | | 4.1 U | 0.082 U | 0.082 U | 19 U | 0.33 U | 0.33 U | 0.082 U | 0.082 U | 3.3 U | 0.082 U | 3.3 U | - | 0.082 U | | | |
| Ethanol | NA | NA | | 15 | 16 | 16 | 29.2 | 23.9 | 88.4 | 35 | 41.8 | 30 | 21.7 | 10 U | 33.7 | 37.1 | 4.7 | 85 | 34.3 J | 81.4 E | 48.6 | 93.5 | 19.2 | | | |
| Ethylbenzene | NA | NA | | 20 | 10 | 15 | 0.48 J | 0.74 U | 0.74 U | 0.56 J | 41.8 U | 1.4 | 0.18 U | 50.4 J | 79.5 | 73.4 | 4.7 0.18 U | 12 | 7.4 U | 10 | 7.4 U | 13 | 0.18 U | | | |
| Ethyl Acetate | NA | NA | | 0.9 U | 19 | 0.9 U | 7.2 | 5.4 | 12 | 6.8 | 4.8 U | 4.3 | 9.4 | 32 U | 1.1 U | 1.1 U | 0.76 | 5.4 | 7.4 0 | 5 | 30 | 4 | 7.9 | | | |
| 4-Ethyltoluene | NA | NA | | 9.3 | 6.9 | 5.9 | 0.084 U | 0.33 U | 0.33 U | 0.084 U | 2.1 U | 4.3 0.084 U | 0.084 U | 9.8 U | 23 | 1.1 0 | 0.084 U | 3.4 4.1 | 3.3 U | 3.5 | 3.3 U | 4.4 | 0.084 U | | | |
| Freon 113 | NA | NA | | 0.84 U | 0.84 U | 0.84 U | 0.16 U | 0.66 U | 0.66 U | | - | 0.16 U | 0.76 J | 19 U | 0.66 U | 0.66 U | 0.16 U | 0.16 U | 6.6 U | 0.16 U | 6.6 U | | 0.16 U | | | |
| Freon 114 | NA | NA | | 0.7 U | 0.7 U | 0.7 U | 0.22 U | 0.00 U | 0.91 U | | - | 0.10 U | 0.22 U | 26 U | 0.91 U | 0.91 U | 0.22 U | 0.10 U | 9.1 U | 0.10 U | 9.1 U | 0.22 U | 0.22 U | | | |
| Heptane | NA | NA | | 19 | 19 | 15 | 0.49 J | 0.31 U | 0.33 U | 0.22 U | 1480 | 1.4 | 0.41 J | 101 | 63.9 | 61.9 | 0.082 U | 9.8 | 3.1 U | 8.6 | 27 J | 9.8 | 0.082 U | | | |
| Hexachlorobutadiene | NA | NA | | 1.4 U | 1.4 U | 1.4 U | 0.21 U | | 0.86 U | | | 0.21 U | 0.21 U | 26 U | 0.86 U | 0.86 U | 0.21 U | 0.21 U | 8.6 U | 0.21 U | 8.6 U | | 0.21 U | | | |



<u>Table 3-1</u> Soil Vapor and Indoor Air Sampling Results Summary

| | Historical Regula | | 2013 and 2015 NYSDOH Revised Standards | Initia | al screening san (sub-slab only) | | | March 2016 sampling event including sub-slab and IAQ samples | | | | | | | | | April 2017 sampling event including sub-slab and IAQ samples | | | | | | | | |
|---------------------------------|--------------------------------|----------------------------|---|------------------|-------------------------------------|------------------|-----------------|--|-----------------|------------------|-----------------|------------------|-------------------|------------------|------------------|------------------|--|------------------|-----------------|------------------|-----------------|------------------|--------------------|--|--|
| Analyte Method TO-15 (ug/m³) | Matrix 1 & 2 Sub-Slab Vapor | Matrix 1 & 2 Indoor Air | Air | SV-6 10/30/15 | SV-7 10/30/15 | SV-8 10/30/15 | SV-9* 3/4/16 | IAQ-9* 3/4/16 | SV-10 3/9/16 | IAQ-10 3/9/16 | SV-11 3/4/16 | IAQ-11 3/4/16 | Ambient 3/9/16 | SV-12 3/17/16 | SV-13 3/17/16 | SV-14 3/17/16 | Ambient 3/17/16 | IAQ-3 4/14/17 | SV-4 4/14/17 | IAQ-4 4/14/17 | SV-5 4/14/17 | IAQ-5 4/14/17 | Ambient 4/14/17 | | |
| Hexane | NA | NA | | 5.6 | 7.4 | 6.7 | 0.92 | 3 | 2 J | 2.4 | 5780 | 2.9 | 1.1 | 1720 | 20 | 22 | 0.67 J | 13 | 3.2 U | 12 | 16 J | 13 | 0.6 J | | |
| 2-Hexanone | NA | NA | | 27 | 20 | 14 | 2.5 | 0.74 U | 0.74 U | 0.18 U | 4.5 U | 0.18 U | 0.18 U | 22 U | 2.5 J | 9 | 0.18 U | 0.18 U | 7.4 U | 0.18 U | 7.4 U | 11 | 0.18 U | | |
| Isopropyl Alcohol | NA | NA | | 1.2 U | 1.2 U | 3.9 | 20 | 34.2 | 62.2 | 4.4 | 9.8 U | 18 | 4.9 | 44 U | 5.2 | 9.3 | 0.81 | 18 | 76 | 17 | 83.3 | 21 | 1.9 | | |
| Methylene chloride | NA | NA | | 1.9 U | 1.9 U | 1.9 U | 0.63 J | 0.35 U | 1.5 J | 5.6 | 2.2 U | 0.76 | 1.7 | 10 U | 0.35 U | 0.35 U | 1.4 | 2.6 | 3.5 U | 2.7 | 3.5 U | 2.2 | 0.76 | | |
| Methyl ethyl ketone | NA | NA | | 56.9 | 64 | 36.9 | 392 | 0.56 U | 776 | 2.2 | 1010 | 0.77 | 1.3 | 17 U | 23 | 36.3 | 0.14 U | 1.2 | 5.6 U | 1.2 | 47.8 | 3.2 | 0.56 J | | |
| Methyl Isobutyl Ketone | NA | NA | | 0.45 U | 0.45 U | 0.45 U | 4.9 | 0.9 U | 4.9 | 0.23 U | 5.7 U | 0.23 U | 0.23 U | 27 U | 8.6 | 4.9 | 0.23 U | 0.23 U | 9 U | 0.23 U | 9 U | 7.8 | 0.23 U | | |
| Methyl Tert Butyl Ether | NA | NA | | 0.36 U | 0.36 U | 0.36 U | 0.072 U | 0.28 U | 0.28 U | 0.072 U | 1.8 U | 0.072 U | 0.072 U | 8.3 U | 0.28 U | 0.28 U | 0.072 U | 0.072 U | 2.8 U | 0.072 U | 2.8 U | 0.072 U | 0.072 U | | |
| Methylmethacrylate | NA | NA | | 0.49 U | 0.49 U | 0.49 U | 0.16 U | 0.66 U | 0.66 U | 0.16 U | 4.1 U | 0.16 U | 0.16 U | 19 U | 0.66 U | 0.66 U | 0.16 U | 0.16 U | 6.6 U | 0.16 U | 6.6 U | 1 | 0.16 U | | |
| Propylene | NA | NA | | 0.55 U | 4 | 22.2 | 2.6 | 7.9 | 4.5 | 5.5 | 1.4 U | 0.055 U | 0.055 U | 6.5 U | 5.5 | 12 | 0.055 U | 36.8 | 2.2 U | 33.3 | 23 J | 39 | 1.1 | | |
| Styrene | NA | NA | | 0.43 U | 0.43 U | 0.43 U | 0.064 U | 0.26 U | 0.26 U | 0.064 U | 1.7 U | 0.064 U | 0.064 U | 7.7 U | 0.26 U | 0.26 U | 0.064 U | 2 | 2.6 U | 1.7 | 2.6 U | 2.9 | 0.064 U | | |
| 1,1,1-Trichloroethane | NA | NA | | 0.71 U | 0.71 U | 0.71 U | 0.13 U | 0.51 U | 0.51 U | 0.13 U | 3.3 U | 0.13 U | 0.13 U | 15 U | 0.51 U | 0.51 U | 0.13 U | 0.13 U | 5.1 U | 0.13 U | 5.1 U | 0.13 U | 0.13 U | | |
| 1,1,2,2-Tetrachloroethane | NA | NA | | 0.82 U | 0.82 U | 0.82 U | 0.11 U | 0.44 U | 0.44 U | 0.11 U | 2.8 U | 0.11 U | 0.11 U | 13 U | 0.44 U | 0.44 U | 0.11 U | 0.11 U | 4.4 U | 0.11 U | 4.4 U | 0.96 J | 0.11 U | | |
| 1,1,2-Trichloroethane | NA | NA | | 0.76 U | 0.76 U | 0.76 U | 0.21 U | 0.87 U | 0.87 U | 0.21 U | 5.4 U | 0.21 U | 0.21 U | 25 U | 0.87 U | 0.87 U | 0.21 U | 0.21 U | 8.7 U | 0.21 U | 8.7 U | 0.21 U | 0.21 U | | |
| 1,2,4-Trichlorobenzene | NA | NA | | 1.3 U | 1.3 U | 1.3 U | 0.42 U | 1.6 U | 1.6 U | 0.42 U | 10 U | 0.42 U | 0.42 U | 49 U | 1.6 U | 1.6 U | 0.42 U | 0.42 U | 16 U | 0.42 U | 16 U | 2.2 | 0.42 U | | |
| 1,2,4-Trimethylbenzene | NA | NA | | 29 | 21 | 17 | 1.6 | 0.3 U | 0.3 U | 0.074 U | 1.9 U | 1.6 | 0.074 U | 8.8 U | 80.1 | 62.9 | 0.074 U | 13 | 3 U | 11 | 3 U | 12 | 0.074 U | | |
| 1,3,5-Trimethylbenzene | NA | NA | | 9.3 | 6.9 | 5.9 | 0.48 J | 0.88 U | 0.88 U | 0.22 U | 5.4 U | 0.54 J | 0.22 U | 26 U | 24 | 19 | 0.22 U | 4 | 8.8 U | 3.4 | 8.8 U | 4.3 | 0.22 U | | |
| 2,2,4-Trimethylpentane | NA | NA | | 2.7 J | 2.6 J | 3.1 J | 0.11 U | 2.2 J | 0.43 U | 2.3 | 2.7 U | 5.1 | 0.84 J | 13 U | 13 | 11 | 0.11 U | 23 | 4.3 U | 20 | 4.3 U | 23 | 0.51 J | | |
| Tertiary Butyl Alcohol | NA | NA | | 22 | 29 | 25 | 20 | 0.64 U | 21 | 0.16 U | 56.4 | 0.16 U | 0.16 U | 19 U | 3.6 | 7.6 | 0.16 U | 0.16 U | 6.4 U | 0.16 U | 6.4 U | 1.4 | 0.16 U | | |
| Tetrachloroethylene | >1000 Mitigate | >100 Mitigate | 30 | 753 | 16 | 2620 | 1460 | 1.1 | 1500 | 0.75 | 7.5 | 0.49 | 0.95 | 18 U | 14 | 258 | 0.16 U | 9.5 | 8070 | 8.8 | 6520 | 8.1 | 1.2 | | |
| Tetrahydrofuran | NA | NA | | 0.5 U | 0.5 U | 0.5 U | 1170 | 0.53 U | 1050 | 1.5 | 2450 | 0.13 U | 0.13 U | 16 U | 0.53 U | 0.53 U | 0.13 U | 0.13 U | 5.3 U | 0.13 U | 5.3 U | 0.86 | 0.13 U | | |
| Toluene | NA | NA | | 69.3 | 67.5 | 58.4 | 3.4 | 7.9 | 6 | 4.9 | 1.2 U | 12 | 2.8 | 199 | 357 | 344 | 0.49 J | 63.3 | 1.9 U | 53.9 | 15 J | 62.9 | 1.6 | | |
| Trichloroethylene | >250 Mitigate | >5 Mitigate | 2 | 1.3 | 0.54 U | 157 | 1 | 0.4 U | 4.3 | 0.1 U | 10 | 0.1 U | 0.1 U | 12 U | 0.4 U | 7.5 | 0.1 U | 0.1 U | 107 | 0.13 J | 4 U | 0.39 | 0.1 U | | |
| Trichlorofluoromethane | NA | NA | | 1.8 J | 0.46 U | 0.46 U | 1.5 | 0.49 U | 0.49 U | 1.9 | 3.1 U | 1.4 | 2.1 | 15 U | 0.49 U | 0.49 U | 0.12 U | 1.3 | 4.9 U | 1.3 | 4.9 U | 1.3 | 1.3 | | |
| Vinyl chloride | >250 Mitigate | >5 Mitigate | | 0.33 U | 0.33 U | 0.33 U | 0.054 U | 0.21 U | 0.21 U | 0.054 U | 23 | 0.054 U | 0.054 U | 854 | 0.21 U | 0.21 U | 0.054 U | 0.054 U | 2.1 U | 0.054 U | 2.1 U | 0.054 U | 0.054 U | | |
| Vinyl Acetate | NA | NA | | 0.77 U | 0.77 U | 0.77 U | 0.19 U | 0.77 U | 0.77 U | 0.19 U | 4.9 U | 0.19 U | 0.19 U | 23 U | 0.77 U | 0.77 U | 0.19 U | 0.19 U | 7.7 U | 0.19 U | 7.7 U | 0.19 U | 0.19 U | | |
| m,p-Xylene | NA | NA | | 80.8 | 69.1 | 57.8 | 2 | 3.5 | 1.9 J | 1.7 | 7.4 U | 6.1 | 0.96 | 182 | 297 | 264 | 0.3 U | 41 | 12 U | 35 | 12 U | 42 | 0.96 | | |
| o-Xylene | NA | NA | | 27 | 24 | 20 | 1.3 | 0.87 U | 0.87 U | 0.65 J | 5.6 U | 2.1 | 0.22 U | 54.3 J | 87.3 | 76.4 | 0.22 U | 15 | 8.7 U | 13 | 8.7 U | 16 | 0.22 U | | |
| Xylenes (total) | NA | NA | | 108 | 93.4 | 77.7 | 3.2 | 3.5 | 1.9 J | 2.3 | 5.6 U | 8.3 | 0.96 | 236 | 384 | 340 | 0.22 U | 56 | 8.7 U | 47.8 | 8.7 U | 58.2 | 0.96 | | |

Notes
* - Sub-slab and IAQ results transposed in lab; values tabulated correspond to what is believed to be the correct values
Sample locations identified in light gray shaded headings are sub-slab soil vapor sample locations.
Sample locations identified in dark gray shaded headings are indoor air quality or ambient air sample locations.
In accordance with NYSDOH letter dated 9/12/2017, highlights of data for sub-slab vapors exceeding historic mitigation guidance levels have been removed.

All results recorded in ug/m³



<u> Table 3-2</u> Soil Analytical Results Summary

| Analyte | NY SCO - Residential w/CP-51 (10/10) | NY SCO - Commercial w/CP-51 (10/10) | NY SCO - Protection of Groundwater | | Initial | Site Screening | g Data (Octobe | r 2015) | - | | _ | January 201 | 6 Investigation | | |
|------------------------------|--|---|---|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|----------------------------|-----------------------------|-----------------------------|-----------------------------|-------------------------------|-----------------------------|-----------------------------|
| VOCs | (6 NYCRR 375-6 12/06) | (6 NYCRR 375-6 12/06) | w/CP-51 (10/10) (6 NYCRR 375-6 12/06) | SB-1 1 ft 10/29/2015 | SB-1 5 ft 10/29/2015 | SB-2 1 ft 10/29/2015 | SB-2 5 ft 10/29/2015 | SB-3 1 ft 10/29/2015 | SB-3 5 ft 10/29/2015 | MW-1I 20 ft 1/18/2016 | MW-1I 45 ft 1/18/2016 | MW-1I 50 ft 1/18/2016 | MW-4S 23.5 ft 1/19/2016 | MW-4S 30 ft 1/19/2016 | MW-4S 40 ft 1/19/2016 |
| Acetone | 100 | 500 | 0.05 | 0.0362 | 0.0025 U | 0.0751 | 0.0196 | 0.014 | 0.0116 | 0.003 U | 0.0027 U | 0.0057 J | 0.023 | 0.0581 | 0.0057 J |
| Benzene | 2.9 | 44 | 0.06 | 0.00014 U | 0.00015 U | 0.00014 U | 0.00014 U | 0.00014 U | 0.00014 U | 0.00018 U | 0.00016 U | 0.00021 U | 0.00018 U | 0.00022 U | 0.00013 U |
| Bromochloromethane | - | - | - | 0.00032 U | 0.00035 U | 0.00033 U | 0.00033 U | 0.00033 U | 0.00032 U | 0.00042 U | 0.00037 U | 0.00049 U | 0.00042 U | 0.0005 U | 0.00029 U |
| Bromodichloromethane | - | - | - | 0.00016 U | 0.00018 U | 0.00016 U | 0.00016 U | 0.00017 U | 0.00016 U | 0.00021 U | 0.00019 U | 0.00025 U | 0.00021 U | 0.00025 U | 0.00015 U |
| Bromoform | - | - | - | 0.00025 U | 0.00027 U | 0.00025 U | 0.00025 U | 0.00025 U | 0.00024 U | 0.00032 U | 0.00028 U | 0.00038 U | 0.00032 U | 0.00038 U | |
| Bromomethane | - | - | - | 0.00038 U | 0.00041 U | 0.00038 U | 0.00038 U | 0.00039 U | 0.00037 U | 0.00049 U | 0.00043 U | 0.00058 U | 0.00049 U | 0.00059 U | 0.00035 U |
| 2-Butanone (MEK) | 100 | 500 | 0.3 | 0.002 U | 0.0022 U | 0.002 U | 0.002 U | 0.0021 U | 0.002 U | 0.0026 U | 0.0023 U | 0.003 U | 0.0026 U | 0.0031 U | 0.0018 U |
| Carbon disulfide | 100 | | 2.7 | 0.00024 U | 0.00026 U | 0.00024 U | 0.00024 U | 0.00024 U | 0.00023 U | 0.00031 U | 0.00027 U | 0.00036 U | 0.00031 U | 0.00037 U | 0.00022 U |
| Carbon tetrachloride | 1.4 | 22 | 0.76 | 0.00024 U | 0.00026 U | 0.00024 U | 0.00024 U | 0.00025 U | | 0.00031 U | 0.00027 U | 0.00037 U | | 0.00037 U | |
| Chlorobenzene | 100 | 500 | 1.1 | 0.00016 U | 0.00018 U | 0.00016 U | 0.00016 U | 0.00017 U | 0.00016 U | 0.00021 U | 0.00018 U | 0.00025 U | | 0.00025 U | 0.00015 U |
| Chloroethane | | | 1.9 | 0.0005 U | 0.00054 U | 0.00051 U | 0.00051 U | 0.00052 U | | 0.00065 U | 0.00057 U | 0.00077 U | | 0.00078 U | |
| Chloroform | 10 | 350 | 0.37 | 0.00016 U | 0.00017 U | 0.00016 U | 0.00016 U | 0.00016 U | 0.00015 U | | 0.00018 U | 0.00024 U | | | 0.00014 U |
| Chloromethane | - | - | - | 0.00027 U | 0.0003 U | 0.00028 U | 0.00028 U | 0.00028 U | 0.00027 U | 0.00035 U | 0.00031 U | 0.00042 U | 0.00035 U | 0.00043 U | 0.00025 U |
| Cyclohexane | - | - | - | 0.00033 U | 0.00036 U | 0.00033 U | 0.00033 U | 0.00034 U | 0.00032 U | 0.00043 U | 0.00038 U | 0.0005 U | | 0.00051 U | |
| 1,2-Dibromo-3-chloropropane | - | - | - | 0.00057 U | 0.00062 U | 0.00058 U | 0.00057 U | 0.00058 U | 0.00056 U | 0.00073 U | 0.00065 U | 0.00087 U | | 0.00089 U | 0.00052 U |
| Dibromochloromethane | | | | 0.00021 U | 0.00023 U | 0.00022 U | 0.00022 U | 0.00022 U | 0.00021 U | 0.00028 U | 0.00024 U | 0.00033 U | 0.00028 U | 0.00033 U | 0.00019 U |
| 1,2-Dibromoethane | - | - | - | 0.00014 U | 0.00015 U | 0.00014 U | 0.00014 U | 0.00014 U | 0.00013 U | 0.00018 U | 0.00016 U | 0.00021 U | | 0.00021 U | 0.00012 U |
| 1,2-Dichlorobenzene | 100 | 500 | 1.1 | 0.00013 U | 0.00014 U | 0.00013 U | 0.00013 U | 0.00013 U | 0.00013 U | 0.00016 U | 0.00015 U | 0.00019 U | 0.00016 U | 0.0002 U | |
| 1,3-Dichlorobenzene | 17 | 280 | 2.4 | 0.00016 U | 0.00018 U | 0.00017 U | 0.00017 U | 0.00017 U | 0.00016 U | 0.00021 U | 0.00019 U | 0.00025 U | | 0.00026 U | 0.00015 U |
| 1,4-Dichlorobenzene | 9.8 | 130 | 1.8 | 0.00024 U | 0.00025 U | 0.00024 U | 0.00024 U | 0.00024 U | 0.00023 U | 0.0003 U | 0.00027 U | 0.00036 U | | 0.00037 U | 0.00021 U |
| Dichlorodifluoromethane | - | - | - | 0.00038 U | 0.00041 U | 0.00038 U | 0.00038 U | 0.00039 U | | 0.00049 U | 0.00043 U | 0.00058 U | 0.00049 U | 0.00059 U | |
| 1,1-Dichloroethane | 19 | 240 | 0.27 | 0.00015 U | 0.00016 U | 0.00015 U | 0.00015 U | 0.00015 U | 0.0001 | | | 0.00023 U | | | 0.00013 U |
| 1,2-Dichloroethane | 2.3 | 30 | 0.02 | 0.00014 U | 0.00015 U | 0.00014 U | 0.00014 U | 0.00014 U | 0.00014 U | 0.00018 U | 0.00016 U | 0.00021 U | 0.00018 U | 0.00022 U | |
| 1,1-Dichloroethene | 100 | 500 | 0.33 | 0.00062 U | 0.00067 U | 0.00063 U | 0.00063 U | 0.00064 U | | 0.0008 U | 0.0007 U | 0.00095 U | 0.0008 U | 0.00096 U | |
| cis-1,2-Dichloroethene | 59 | 500 | 0.25 | 0.00082 U | 0.00088 U | 0.00082 U | 0.00082 U | 0.00084 U | 0.0000 | | 0.00093 U | 0.0012 U | 0.0011 U | 0.0013 U | |
| trans-1,2-Dichloroethene | 100 | 500 | 0.19 | 0.00062 U | 0.00067 U | 0.00063 U | 0.00063 U | 0.00064 U | 0.00061 U | 0.0008 U | 0.00071 U | 0.00095 U | | 0.00097 U | |
| 1,2-Dichloropropane | | | | 0.00025 U | 0.00027 U | 0.00025 U | 0.00025 U | 0.00026 U | 0.00024 U | 0.00032 U | 0.00028 U | 0.00038 U | | | 0.00023 U |
| cis-1,3-Dichloropropene | - | - | - | 0.00012 U | 0.00013 U | 0.00012 U | 0.00012 U | 0.00013 U | | 0.00016 U | 0.00014 U | 0.00019 U | 0.00016 U | 0.00019 U | |
| trans-1,3-Dichloropropene | - | - | - | 0.00019 U | 0.0002 U | 0.00019 U | 0.00019 U | 0.00019 U | 0.00018 U | 0.00024 U | 0.00021 U | 0.00028 U | | 0.00029 U | |
| Ethylbenzene | 30 | 390 | 1 | 0.00017 U | 0.00018 U | 0.00017 U | 0.00017 U | 0.00018 U | 0.00017 U | 0.00022 U | 0.00019 U | 0.00026 U | 0.00022 U | 0.00027 U | |
| Freon 113 | 100 | | 6 | 0.00047 U | 0.00051 U | 0.00047 U | 0.00047 U | 0.00048 U | | 0.0006 U | 0.00053 U | 0.00072 U | 0.00061 U | 0.00073 U | |
| 2-Hexanone | - | - | - | 0.0014 U | 0.0015 U | 0.0014 U | 0.0014 U | 0.0014 U | | 0.0018 U | | 0.0021 U | 0.0018 U | 0.0022 U | |
| Isopropylbenzene | 100 | | 2.3 | 0.00011 U | 0.00012 U | 0.00011 U | 0.00011 U | 0.00011 U | 0.00011 U | 0.00014 U | 0.00013 U | 0.00017 U | 0.00014 U | 0.00017 U | 0.0001 U |
| Methyl Acetate | - | - | - | 0.0009 U | 0.00097 U | 0.00091 U | 0.00091 U | 0.00092 U | 0.00088 U | 0.0012 U | 0.001 U | 0.0014 U | 0.0012 U | 0.0014 U | |
| Methylcyclohexane | - | - | - | 0.00024 U | 0.00026 U | | 0.00024 U | 0.00024 U | | | | | | | 0.00022 U |
| Methyl Tert Butyl Ether | 62 | 500 | 0.93 | | | | | | | | | | 0.00021 U | | |
| 4-Methyl-2-pentanone(MIBK) | E4 | 500 | 1 | | | | | | | | | | 0.00062 U | | |
| Methylene chloride | 51 | 500 | 0.05 | 0.001 U 0.00019 U | 0.0011 U | 0.001 U 0.00019 U | | 0.0011 U 0.00019 U | | 0.0013 U | | 0.0016 U | 0.00013 U 0.00024 U | | 0.00093 U |
| Styrene | 25 | | 0.6 | | | | | 0.00019 U | | | | | | | |
| 1,1,2,2-Tetrachloroethane | 35 5.5 | 150 | 0.6 | 0.00018 U 0.0183 | 0.0002 U | 0.00018 U 0.0347 | 0.00018 0 | | | | 0.00021 U 0.0043 | | 0.00024 U 0.00041 U | | |
| Tetrachloroethene Toluene | 5.5 100 | 500 | 1.3 0.7 | | 0.0196 0.00024 U | | | 0.0233 | 0.0238 | 0.0139 | | | 0.00041 U 0.00028 U | | |
| 1,2,3-Trichlorobenzene | 100 | 500 | 0.7 | 0.00022 U 0.00018 U | 0.00024 U | | 0.00022 U 0.00019 U | 0.00022 U 0.00019 U | | | | | 0.00028 U 0.00024 U | | 0.0002 U 0.00017 U |
| 1,2,3-Trichlorobenzene | + | | 3.4 | 0.00018 U | 0.0002 U 0.00019 U | | 0.00019 U 0.00018 U | | 0.00018 U 0.00017 U | 0.00024 U 0.00023 U | | | | | 0.00017 U 0.00016 U |
| 1,1,1-Trichloroethane | 100 | 500 | 0.68 | | | 0.00018 U | | | | | | | 0.00023 U | | |
| 1,1,2-Trichloroethane | - | - 500 | 0.00 | | 0.00017 U | 0.00016 U | | | 0.00015 U | | 0.00018 U 0.00017 U | | | | 0.00014 U |
| Trichloroethene | - 10 | 200 | - 0.47 | 0.00015 U | 0.00017 U 0.0004 J | | 0.00018 U | | 0.00015 U 0.00039 J | 0.0002 U 0.00074 J | | 0.00023 U | | | |
| Trichlorofluoromethane | | | | | 0.0004 J 0.00028 U | | 0.00088 J | 0.00016 U 0.00027 U | | | | | 0.0002 0 | | 0.00014 U |
| Vinyl chloride | - 0.21 | - 13 | - 0.02 | | | | | | | 0.00034 U 0.00027 U | | | 0.00034 U 0.00027 U | | |
| | 100 | 500 | 1.6 | 0.00021 U 0.00029 U | | 0.00021 U 0.00029 U | | 0.00021 U 0.00029 U | | | | | 0.00027 U 0.00037 U | | |
| Xylene (total) | 100 | 500 | 0.1 | 0.00029 U | 0.00031 U | 0.00029 U | 0.00029 0 | 0.00029 U | 0.00028 U | 0.00037 U | 0.00033 0 | 0.00044 U | 0.00037 U | 0.00045 U | 0.00020 U |

NOTE:

All values recorded in units of mg/kg

U - Value reported under the detection limit

J - Approximated value

Cells highlighted in green and italics indicate an exceedance of the parameter's Protection of Groundwater SCO, but not the Residential or Commercial SCO. Cells highlighted in yellow and bold indicate exceedance of the parameter's Protection of Groundwater SCO and the Residential SCO, but not the Commercial SCO.

Cells highlighted in light red and bold italics indicate an exceedance of the parameter's Protection of Groundwater SCO, Residential SCO, and Commercial SCO.



<u> Table 3-2</u> Soil Analytical Results Summary

| Analyte | NY SCO - Residential w/CP-51 (10/10) | NY SCO - Commercial w/CP-51 (10/10) | NY SCO - Protection of Groundwater | | | | Mar | ch 2016 Investi | gation | | |
|-----------------------------|--|---|---|-----------------------------|-------------------|-----|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|
| VOCs | (6 NYCRR 375-6 12/06) | (6 NYCRR 375-6 12/06) | w/CP-51 (10/10) (6 NYCRR 375-6 12/06) | MW-5 22.5 ft 3/8/2016 | MV 30 3/8/2 | ft | MW-5 35 ft 3/8/2016 | MW-6 23 ft 3/8/2016 | MW-6 30 ft 3/8/2016 | MW-7 23 ft 3/8/2016 | MW-7 34 ft 3/8/2016 |
| Acetone | 100 | 500 | 0.05 | 0.0129 J | 0.002 | | 0.0029 U | 0.203 | 0.283 | 0.029 U | 0.033 U |
| Benzene | 2.9 | 44 | 0.06 | 0.00017 U | 0.0001 | | 0.00017 U | 0.0022 U | 0.003 U | 0.0017 U | 0.002 U |
| Bromochloromethane | - | | - | 0.0004 U | 0.0004 | | 0.0004 U | 0.0052 U | 0.0069 U | 0.004 U | 0.0045 U |
| Bromodichloromethane | | - | - | 0.0004 U | 0.000 | - | 0.0004 U | 0.0032 U | 0.0035 U | 0.004 U | 0.0023 U |
| Bromoform | | - | - | 0.0002 U | 0.0003 | | 0.00031 U | 0.004 U | 0.0053 U | 0.002 U | 0.0035 U |
| Bromomethane | - | | | 0.00048 U | 0.0004 | | 0.00048 U | 0.0061 U | 0.0082 U | 0.0047 U | 0.0053 U |
| 2-Butanone (MEK) | 100 | 500 | 0.3 | 0.0025 U | 0.0004 | | 0.0025 U | 0.032 U | 0.0032 0 | 0.0047 U | 0.028 U |
| Carbon disulfide | 100 | 500 | 2.7 | 0.00023 U | 0.002 | | 0.0023 U | 0.0032 U | 0.0051 U | 0.0029 U | 0.0034 U |
| Carbon tetrachloride | 1.4 | 22 | 0.76 | 0.0003 U | 0.000 | | 0.0003 U | 0.0039 U | 0.0051 U | 0.0029 U | 0.0034 U |
| Chlorobenzene | 1.4 | 500 | 1.1 | 0.0003 U | 0.000 | | 0.0003 U | 0.0039 U | 0.0032 U | 0.0029 U | 0.0023 U |
| Chloroethane | 100 | 500 | 1.9 | 0.00063 U | 0.0006 | - | 0.0002 0 | 0.0020 U | 0.0033 0 | 0.002 U | 0.0023 0 0.0071 U |
| Chloroform | 10 | 350 | 0.37 | 0.00003 U 0.00019 U | 0.0001 | | 0.00019 U | 0.0025 U | 0.0033 U | 0.0002 U | 0.0022 U |
| Chloromethane | | | | 0.00019 U 0.00034 U | 0.0003 | | 0.00034 U | 0.0023 U 0.0044 U | 0.0059 U | 0.0019 U | 0.0022 U |
| Chloromethane | - | - | - | 0.00034 U 0.00041 U | 0.0003 | | 0.00034 U 0.00041 U | 0.0044 U 0.0053 U | 0.0059 U 0.0071 U | 0.0034 U 0.004 U | 0.0039 U 0.0046 U |
| 1,2-Dibromo-3-chloropropane | - | - | - | 0.00041 U 0.00071 U | 0.0004 | | 0.00041 U | 0.0053 U 0.0091 U | 0.0071 U | 0.004 U 0.007 U | 0.0046 U 0.008 U |
| | - | - | - | | | - | | | | | |
| Dibromochloromethane | | | | 0.0002. 0 | 0.0002 | | ••••• | 0.0034 U 0.0022 U | | | |
| 1,2-Dibromoethane | - 100 | - | - | 0.0001.0 | 0.0001 | | 0.00017 U | | 0.0029 U 0.0027 U | 0.0017 U | |
| 1,2-Dichlorobenzene | | 500 | 1.1 | 0.00016 U | 0.0001 | | 0.00031 J | 0.002 U | 010021 0 | 0.0016 U | 0.0018 U |
| 1,3-Dichlorobenzene | 17 | 280 | 2.4 | 0.00021 U | 0.0002 | | 0.00021 U | 0.0026 U | 0.0035 U | 0.002 U | 0.0023 U |
| 1,4-Dichlorobenzene | 9.8 | 130 | 1.8 | 0.00029 U | 0.0002 | | 0.00029 U | 0.0038 U | 0.005 U | 0.0029 U | 0.0033 U |
| Dichlorodifluoromethane | - | - | - | 0.00047 U | 0.0004 | | 0.00047 U | 0.0061 U | 0.0081 U | 0.0046 U | 0.0053 U |
| 1,1-Dichloroethane | 19 | 240 | 0.27 | 0.00018 U | 0.0001 | | 0.00018 U | 0.0024 U | 0.0032 U | 0.0018 U | 0.0021 U |
| 1,2-Dichloroethane | 2.3 | 30 | 0.02 | 0.00018 U | 0.0001 | | 0.00018 U | 0.0023 U | 0.003 U | 0.0017 U | 0.002 U |
| 1,1-Dichloroethene | 100 | 500 | 0.33 | 0.00077 U | 0.0007 | 7 U | 0.0025 | 0.0099 U | 0.013 U | 0.0076 U | 0.0087 U |
| cis-1,2-Dichloroethene | 59 | 500 | 0.25 | 0.001 U | 2.44 | | 3.67 | 0.013 U | 0.017 U | 0.01 U | 0.011 U |
| trans-1,2-Dichloroethene | 100 | 500 | 0.19 | 0.00078 U | 0.0007 | | 0.0044 | 0.01 U | 0.013 U | 0.0076 U | 0.0087 U |
| 1,2-Dichloropropane | | | | 0.00031 U | 0.0003 | | 0.00031 U | 0.004 U | 0.0053 U | 0.003 U | 0.0035 U |
| cis-1,3-Dichloropropene | - | - | - | 0.00015 U | 0.0001 | | 0.00015 U | 0.002 U | 0.0026 U | 0.0015 U | 0.0017 U |
| trans-1,3-Dichloropropene | - | - | - | 0.00023 U | 0.0002 | | 0.00023 U | 0.003 U | 0.004 U | 0.0023 U | 0.0026 U |
| Ethylbenzene | 30 | 390 | 1 | 0.0053 | 0.0002 | | 0.00021 U | 0.01 J | 0.0088 J | 0.0021 U | 0.0024 U |
| Freon 113 | 100 | | 6 | 0.00059 U | 0.0005 | | 0.00059 U | 0.0075 U | 0.01 U | 0.0057 U | 0.0066 U |
| 2-Hexanone | - | - | - | 0.0018 U | 0.001 | | 0.0018 U | 0.023 U | 0.03 U | 0.017 U | 0.02 U |
| Isopropylbenzene | 100 | | 2.3 | 0.00014 U | 0.0001 | | 0.00015 J | 0.0018 U | 0.0024 U | 0.0014 U | 0.0016 U |
| Methyl Acetate | - | - | - | 0.0011 U | 0.001 | U | 0.0011 U | 0.014 U | 0.019 U | 0.011 U | 0.013 U |
| Methylcyclohexane | - | - | - | 0.0003 U | 0.000 | | 0.0003 U | 0.0038 U | 0.0051 U | 0.0029 U | 0.0033 U |
| Methyl Tert Butyl Ether | 62 | 500 | 0.93 | 0.0002 U | | | | | | 0.002 U | 0.0022 U |
| 4-Methyl-2-pentanone(MIBK) | | | 1 | 0.0006 U | 0.000 | | 0.0006 U | 0.0077 U | 0.01 U | 0.0059 U | 0.0067 U |
| Methylene chloride | 51 | 500 | 0.05 | 0.0013 U | 0.001 | | 0.0013 U | 0.016 U | 0.022 U | 0.013 U | 0.014 U |
| Styrene | | | | 0.00023 U | 0.0002 | | 0.00023 U | 0.003 U | 0.004 U | 0.0023 U | 0.0026 U |
| 1,1,2,2-Tetrachloroethane | 35 | | 0.6 | 0.00023 U | 0.0002 | 3 U | 0.00023 U | 0.0029 U | 0.0039 U | 0.0022 U | 0.0026 U |
| Tetrachloroethene | 5.5 | 150 | 1.3 | 0.0114 | 2.74 | | 3.35 | 0.0051 U | 0.0067 U | 0.0039 U | 0.0076 J |
| Toluene | 100 | 500 | 0.7 | 0.117 | 0.0002 | | 0.00027 U | 0.143 | 0.124 | 0.0027 U | 0.0031 U |
| 1,2,3-Trichlorobenzene | | | | 0.00023 U | 0.0002 | | 0.00023 U | 0.003 U | 0.0039 U | 0.0023 U | 0.0026 U |
| 1,2,4-Trichlorobenzene | | | 3.4 | 0.00022 U | 0.0002 | | 0.00022 U | 0.0028 U | 0.0038 U | 0.0022 U | 0.0025 U |
| 1,1,1-Trichloroethane | 100 | 500 | 0.68 | 0.00019 U | 0.0001 | | 0.00019 U | 0.0025 U | 0.0033 U | 0.0019 U | 0.0022 U |
| 1,1,2-Trichloroethane | - | - | - | 0.00019 U | 0.0001 | 9 U | 0.00019 U | 0.0025 U | 0.0033 U | 0.0019 U | 0.0022 U |
| Trichloroethene | 10 | 200 | 0.47 | 0.00019 U | 0.101 | | 0.636 | 0.0025 U | 0.0033 U | 0.0019 U | 0.0022 U |
| Trichlorofluoromethane | - | - | - | 0.00033 U | 0.0003 | 3 U | 0.00033 U | 0.0042 U | 0.0056 U | 0.0032 U | 0.0037 U |
| Vinyl chloride | 0.21 | 13 | 0.02 | 0.00026 U | 0.0002 | | 0.00026 U | 0.0033 U | 0.0044 U | 0.0025 U | 0.0029 U |
| Xylene (total) | 100 | 500 | 1.6 | 0.0265 | 0.0003 | | 0.00036 U | 0.0564 | 0.0436 | 0.0035 U | 0.004 U |

NOTE: All values recorded in units of mg/kg U - Value reported under the detection limit J - Approximated value Cells highlighted in green and italics indicate an exceedance of the parameter's Protection of Groundwater SC Cells highlighted in yellow and bold indicate exceedance of the parameter's Protection of Groundwater SC Cells highlighted in light red and bold italics indicate an exceedance of the parameter's Protection of Groundwater SC



<u> Table 3-2</u> Soil Analytical Results Summary

| Analyte | NY SCO - Residential w/CP-51 (10/10) | NY SCO - Commercial w/CP-51 (10/10) | NY SCO - Protection of Groundwater | | | | | | | | | June 20 | 016 I | nvestigatio | n | | | | | | | | |
|-----------------------------|--|---|---|-----------------------------|---|----------------------------|----------|--------------------------|--------------|-----------------------------|---|------------------------------|-------|----------------------------|---|---|------|---------------------------|---|-------------------------------|----------|-----------------------------|---|
| VOCs | (6 NYCRR 375-6 12/06) | (6 NYCRR 375-6 12/06) | w/CP-51 (10/10) (6 NYCRR 375-6 12/06) | MW-5I 20 ft 6/22/2016 | | MW-5I 23 ft 6/22/201 | | MW-5I ft 6/22/2016 | 6 | MW-5I 50 ft 6/22/2016 | | MW-8 19.5 ft 6/22/2016 | 6 | MW-8 30 ft 5/22/2016 | 6 | MW-9 19.4 ft /22/2016 | | MW-9 30 ft /22/2016 | | MW-10 19.5 ft 6/22/2016 | | MW-10 30 ft 6/22/2016 | 6 |
| Acetone | 100 | 500 | 0.05 | 0.0022 L | J | 0.0022 | U | | U | 0.0026 U | | 0.0023 U | | 0296 | | 0077 J | 0.0 | | | 0.0027 U | 1 | 0.0025 | U |
| Benzene | 2.9 | 44 | 0.06 | 0.00014 | | 0.00014 | - | | U | 0.00017 U | _ | 0.00015 U | | 00013 U | | 0015 U | 0.00 | | | 0.00017 L | - | 0.00016 | U |
| Bromochloromethane | - | - | - | 0.00038 | | 0.00038 | - | 0.00041 | U | 0.00044 U | _ | 0.0004 U | | 00036 U | | 0004 U | 0.00 | | | |] | 0.00042 | U |
| Bromodichloromethane | - | - | - | 0.00018 | | 0.00018 | U | | Ŭ | 0.00021 U | | 0.00019 U | | 00017 U | | 00019 U | 0.00 | | | | J | 0.0002 | Ŭ |
| Bromoform | - | - | - | 0.00032 | | 0.00032 | _ | | Ŭ | 0.00037 U | | 0.00033 U | | 0003 U | | 00033 U | 0.00 | | | 0.00039 | J | 0.00035 | Ŭ |
| Bromomethane | - | - | - | 0.00058 L | - | 0.00058 | - | | Ŭ | 0.00067 U | - | 0.00061 U | - | 00054 U | | 0006 U | 0.00 | | | | J | 0.00064 | Ū |
| 2-Butanone (MEK) | 100 | 500 | 0.3 | 0.0021 | | 0.0021 | - | | Ŭ | 0.0024 U | | 0.0022 U | | .002 U | | 0022 U | 0.0 | | | | J | 0.0023 | Ŭ |
| Carbon disulfide | 100 | | 2.7 | 0.0002 | | 0.0002 | | | U | 0.00023 U | | 0.00021 U | - | 00019 U | | 0021 U | 0.00 | | | | J | 0.00022 | U |
| Carbon tetrachloride | 1.4 | 22 | 0.76 | 0.0002 1 | | 0.0002 | | | Ŭ | 0.00023 U | | 0.00021 U | | 00019 U | | 00021 U | 0.00 | | _ | 0.00024 | - | 0.00022 | Ŭ |
| Chlorobenzene | 100 | 500 | 1.1 | 0.00019 | | 0.00019 | - | 0.00021 | Ŭ | 0.00022 U | | 0.0002 U | - | 00018 U | | 0002 U | 0.00 | | | 0.00024 | J | 0.00021 | Ŭ |
| Chloroethane | | | 1.9 | 0.00051 | | 0.00051 | - | | Ŭ | 0.00059 U | | 0.00054 U | | 00048 U | - | 00053 U | 0.00 | | | | J | 0.00056 | Ŭ |
| Chloroform | 10 | 350 | 0.37 | 0.00028 | | 0.00028 | | | Ŭ | 0.00033 U | _ | 0.0003 U | - | 00027 U | | 0003 U | 0.00 | | _ | | J | 0.00031 | Ŭ |
| Chloromethane | - | - | - | 0.00025 | _ | 0.00025 | | | Ŭ | 0.00029 U | _ | 0.00026 U | - | 00024 U | | 00026 U | 0.00 | | | | J | 0.00028 | U |
| Cyclohexane | - | - | - | 0.00065 | | 0.00065 | | | U | 0.00075 U | _ | 0.00068 U | - | 00061 U | | 00068 U | 0.00 | | | | J | 0.00072 | U |
| 1,2-Dibromo-3-chloropropane | - | - | - | 0.00058 | | 0.00057 | | | Ŭ | 0.00067 U | | 0.00061 U | | 00054 U | | 0006 U | 0.00 | | _ | 0.00071 L | - | 0.00064 | U |
| Dibromochloromethane | | | | 0.00018 | | 0.00018 | - | | Ŭ | 0.00021 U | | 0.00019 U | | 00017 U | - | 00019 U | 0.00 | | _ | | J | 0.0002 | U |
| 1,2-Dibromoethane | - | - | - | 0.00029 | _ | 0.00029 | - | | Ŭ | 0.00033 U | _ | 0.0003 U | | 00027 U | | 0003 U | 0.00 | | | | J | 0.00032 | U |
| 1,2-Dichlorobenzene | 100 | 500 | 1.1 | 0.0002 | | 0.0002 | | | Ŭ | 0.00024 U | _ | 0.00021 U | - | 00019 U | | 00021 U | 0.00 | | | | J | 0.00022 | U |
| 1,3-Dichlorobenzene | 17 | 280 | 2.4 | 0.00016 | | 0.00016 | - | | Ŭ | 0.00019 U | | 0.00017 U | - | 00015 U | | 00017 U | 0.00 | | _ | | J | 0.00018 | U |
| 1,4-Dichlorobenzene | 9.8 | 130 | 1.8 | 0.00018 | | 0.00018 | Ū | 0.0002 | U I | 0.00021 U | | 0.00019 U | | 00017 U | | 00019 U | 0.00 | | | | 1 | 0.0002 | U |
| Dichlorodifluoromethane | - | - | - | 0.00065 | | 0.00065 | <u> </u> | | U | 0.00075 U | | 0.00068 U | | 00061 U | | 00068 U | 0.00 | | | | J | 0.00072 | Ŭ |
| 1,1-Dichloroethane | 19 | 240 | 0.27 | 0.00022 | | 0.00022 | | | U | 0.00026 U | | 0.00023 U | - | 00021 U | | 00023 U | 0.0 | | _ | | J | 0.00025 | Ŭ |
| 1,2-Dichloroethane | 2.3 | 30 | 0.02 | 0.00022 U | | 0.00022 | | | U | 0.00020 U | _ | 0.00020 U | - | 000 <u>21 0</u> 0019 U | | 002 <u>0</u> 0021 U | 0.00 | | | | J | 0.00020 | U |
| 1,1-Dichloroethene | 100 | 500 | 0.33 | 0.00018 L | | 0.00018 | U | | U | 0.00024 U | | 0.00019 U | - | 0001 <u>0</u> 0017 U | | 000 <u>1</u> 9 U | 0.00 | | _ | | J | 0.00022 | U |
| cis-1,2-Dichloroethene | 59 | 500 | 0.25 | 0.0308 | | 0.00052 | - | | U | 0.0006 U | _ | 0.00055 U | - | 00049 U | | 00054 U | 0.00 | | | | J | 0.00058 | Ŭ |
| trans-1,2-Dichloroethene | 100 | 500 | 0.19 | 0.00019 L | | 0.00019 | U | | U | 0.00022 U | _ | 0.0002 U | - | 00018 U | | 0002 U | 0.00 | | | | 1 | 0.00021 | U |
| 1,2-Dichloropropane | 100 | 500 | 0.15 | 0.00037 L | | 0.00037 | U | | U | 0.00043 U | | 0.00039 U | | 00035 U | | 0002 U | 0.00 | | | | , , | 0.00021 | U |
| cis-1,3-Dichloropropene | - | - | - | 0.00023 L | | 0.00023 | - | | U | 0.00043 U | _ | 0.00025 U | - | 00000 U | | 00000 U | 0.00 | | | 0.00029 L | - | 0.00041 | U |
| trans-1,3-Dichloropropene | - | - | - | 0.00026 | | 0.00026 | - | | U | 0.0003 U | | 0.00028 U | | 00025 U | | 002 <u>7</u> U | 0.00 | | _ | | J | 0.00029 | Ŭ |
| Ethylbenzene | 30 | 390 | 1 | 0.00018 L | | 0.00018 | - | | U | 0.00021 U | _ | 0.00019 U | | 000 <u>20 0</u> | | 000 <u>18</u> U | 0.00 | | | | 1 | 0.0002 | Ŭ |
| Freon 113 | 100 | 000 | 6 | 0.00058 | | 0.00057 | - | | U | 0.00067 U | | 0.00061 U | | 00054 U | | 0006 U | 0.00 | | | | J | 0.00064 | U |
| 2-Hexanone | - | - | - | 0.0017 L | _ | 0.0017 | | | U | 0.0019 U | _ | 0.0017 U | - | 0016 U | | 0017 U | 0.0 | | | | J | 0.0018 | Ŭ |
| Isopropylbenzene | 100 | | 2.3 | 0.00018 L | _ | 0.00018 | U | 0.0002 | U I | 0.00021 U | _ | 0.00019 U | - | 0010 U | | 00119 U | 0.00 | | | | <u>,</u> | 0.0002 | U |
| Methyl Acetate | - | - | - | 0.0024 | | 0.0024 | U | | U | 0.0028 U | _ | 0.0025 U | | 0023 U | | 0025 U | 0.0 | | | | 1 | 0.0027 | Ŭ |
| Methylcyclohexane | - | - | - | 0.00024 C | - | 0.00024 | - | | U | 0.00020 U | _ | 0.00063 U | - | 002 <u>5</u> 00057 U | | 002 <u>0</u> 0063 U | 0.0 | | _ | | J | 0.00027 | U |
| Methyl Tert Butyl Ether | 62 | 500 | 0.93 | | | | | | - | 0.00036 U | | 0.00033 U | | 0003 U | | 00033 U | | | _ | | - | 0.00035 | U |
| 4-Methyl-2-pentanone(MIBK) | | | 1 | | | 0.001 | | | Ŭ | | | 0.0011 U | | 00095 U | | 0011 U | | | _ | | J | 0.0011 | Ŭ |
| Methylene chloride | 51 | 500 | 0.05 | 0.0014 | - | 0.0011 | | 0.00098 | J. | 0.0011 J | _ | 0.0012 J | | 00051 J | | 0015 J | 0.0 | | | 0.0012 | ī | 0.0012 | J |
| Styrene | 01 | 000 | 0.00 | 0.00017 U | | | - | | Ŭ | 0.0002 U | | 0.00018 U | | 00016 U | | 0018 U | 0.00 | | | | J | 0.00019 | Ŭ |
| 1,1,2,2-Tetrachloroethane | 35 | | 0.6 | 0.00029 | | | | | U | 0.00033 U | | 0.0003 U | | 00027 U | | 0003 U | 0.00 | | _ | | J | 0.00031 | U |
| Tetrachloroethene | 5.5 | 150 | 1.3 | 0.736 | _ | 0.0036 | 0 | 0.0022 | .1 | 0.0026 J | _ | 0.00035 U | | 000 <u>4</u> 8 J | | 00035 U | 0.0 | | | 0.00077 | <u> </u> | 0.00011 | J |
| Toluene | 100 | 500 | 0.7 | 0.00015 L | | | 11 | 0.00016 | i | 0.00017 U | | 0.00016 U | | 00040 U | | 00000 U | 0.00 | | | | J | 0.00016 | Ŭ |
| 1,2,3-Trichlorobenzene | 100 | 000 | 0.7 | 0.00013 U | | 0.00013 | | | U | 0.00031 U | | 0.00010 U 0.00028 U | | 00014 U | | 001 <u>3</u> 0000000000000000000000000000000000 | 0.00 | | _ | | J | 0.0003 | U |
| 1,2,4-Trichlorobenzene | | | 3.4 | 0.00022 | | | | | - | 0.00025 U | | 0.00023 U | | 0002 <u>3</u> 0 0002 U | | 002 <u>0</u> 0023 U | 0.0 | | | | J | 0.00024 | U |
| 1,1,1-Trichloroethane | 100 | 500 | 0.68 | 0.00022 U | | | | 0.00024 | | | | 0.00023 U | | 000 <u>2</u> 00 00019 U | | 002 <u>3</u> 0000000000000000000000000000000000 | | | | | J | 0.00024 | U |
| 1,1,2-Trichloroethane | - | - | - | 0.00039 L | _ | | | 0.00042 | | 0.00044 U | | 0.0004 U | | 00036 U | | 0004 U | 0.00 | | _ | | J | 0.00042 | U |
| Trichloroethene | - 10 | 200 | 0.47 | 0.00039 | | 0.00036 | | 0.00042 | | 0.00044 U | | 0.00024 U | | 00030 U 00021 U | | 0004 U | 0.00 | | _ | 0.00047 C | _ | 0.00042 | U |
| Trichlorofluoromethane | - | - | - | 0.00039 0.00075 L | _ | | | 0.00023 | | 0.00020 U | | 0.00024 0 0.00079 U | | 00021 U | | 00024 0 00078 U | 0.00 | | _ | | J | 0.00023 | U |
| Vinyl chloride | 0.21 | - 13 | 0.02 | 0.00075 U | _ | | | | | 0.00087 U | _ | 0.00079 U 0.00025 U | | 00071 U | | 0078 U | 0.00 | | _ | | J | 0.00083 | U |
| Xylene (total) | 100 | 500 | 1.6 | 0.00024 U | | 0.00024 | | | | | | 0.00025 U | | 0023 U | | 002 <u>5</u> 0 0025 U | | 0022 U | _ | | J | | U |

NOTE: All values recorded in units of mg/kg U - Value reported under the detection limit J - Approximated value Cells highlighted in green and italics indicate an exceedance of the parameter's Protection of Groundwater SC Cells highlighted in yellow and bold indicate exceedance of the parameter's Protection of Groundwater SC Cells highlighted in light red and bold italics indicate an exceedance of the parameter's Protection of Groundwater SC



| <u>Table 3-3</u> |
|---------------------------------------|
| Soil Sampling Results Summary (SVOCs) |
| |

| Analyte | NY SCO - Residential w/CP-51 (10/10) | NY SCO - Commercial w/CP-51 (10/10) | NY SCO - Industrial w/CP-51 (10/10) | SB-1 | | SB-1 | SB-2 | 2 | SB-2 | 2 | SB-3 | 3 | SB-3 | 3 |
|--|--|---|---|-------------------|--------|----------------------|------------------|--------|------------------|--------|------------------|--------|------------------|--------|
| | (6 NYCRR 375-6 12/06) | (6 NYCRR 375-6 12/06) | (6 NYCRR 375-6 12/06) | 1 ft 10/29/201 | 5 | 5 ft 10/29/2015 | 1 ft 10/29/20 | | 5 ft 10/29/20 | | 1 ft 10/29/20 | | 5 ft 10/29/20 | |
| GC/MS Semi-volatiles (SW846 8270D) (mg/kg) | | | | | | | <u> </u> | | | | | | | |
| 2-Chlorophenol | 100 | | | 0.027 | U | 0.025 U | 0.027 | U | 0.025 | U | 0.027 | U | 0.026 | U |
| 4-Chloro-3-methyl phenol | - | - | - | 0.033 | U | 0.031 U | 0.033 | U | 0.03 | U | 0.033 | U | 0.031 | U |
| 2,4-Dichlorophenol | 100 | | | 0.029 | U | 0.027 U | 0.029 | U | 0.027 | U | 0.029 | U | 0.028 | U |
| 2,4-Dimethylphenol | - | - | - | 0.066 | U | 0.062 U | 0.066 | U | 0.061 | U | 0.066 | U | 0.063 | U |
| 2,4-Dinitrophenol | 100 | | | 0.16 | U | 0.15 U | 0.16 | U | 0.15 | U | 0.16 | U | 0.15 | U |
| 4,6-Dinitro-o-cresol | - | - | - | 0.069 | U | 0.065 U | 0.069 | U | 0.064 | U | 0.069 | U | 0.066 | U |
| 2-Methylphenol | 100 | 500 | 1000 | 0.052 | U | 0.049 U | 0.052 | U | 0.049 | U | 0.052 | U | 0.05 | U |
| 3&4-Methylphenol | - | - | - | 0.034 | U | 0.033 U | 0.035 | U | 0.032 | U | 0.035 | U | 0.033 | U |
| 2-Nitrophenol | | | | 0.033 | U | 0.031 U | 0.033 | U | 0.031 | U | 0.033 | U | 0.032 | U |
| 4-Nitrophenol | | | | 0.061 | U | 0.058 U | 0.061 | U | 0.057 | U | 0.061 | U | 0.059 | U |
| Pentachlorophenol | 2.4 | 6.7 | 55 | 0.088 | U | 0.083 U | 0.088 | U | 0.082 | U | 0.088 | U | 0.085 | U |
| Phenol | 100 | 500 | 1000 | 0.027 | U | 0.025 U | 0.027 | U | 0.025 | U | 0.027 | U | 0.026 | U |
| 2,3,4,6-Tetrachlorophenol | - | - | - | 0.034 | U | 0.032 U | 0.034 | U | 0.032 | U | 0.034 | U | 0.033 | U |
| 2,4,5-Trichlorophenol | 100 | | | 0.032 | U | 0.031 U | 0.033 | U | 0.03 | U | 0.033 | U | 0.031 | U |
| 2,4,6-Trichlorophenol | 100 | | | 0.029 | U | 0.027 U | 0.029 | U | 0.027 | U | 0.029 | U | 0.028 | U |
| Acenaphthene | 100 | 500 | 1000 | 0.020 | U | 0.032 U | 0.034 | U | 0.032 | U | 0.034 | U | 0.033 | U |
| Acenaphthylene | 100 | 500 | 1000 | 0.0034 | U | 0.0032 U | 0.0034 | U | 0.0032 | U | 0.0034 | U | 0.0036 | U |
| Acetophenone | - | - 500 | - | 0.0038 | U | 0.0038 U | 0.0038 | U | 0.0035 | U | 0.0038 | U | 0.0059 | U |
| Acetophenone | - 100 | - 500 | - 1000 | 0.0081 | U | 0.0058 U 0.0029 U | 0.0081 | U | 0.0057 | U | 0.0061 | U | 0.0059 | U |
| Antriacene | | | - | 0.0031 | U | 0.0029 U 0.014 U | 0.0031 | U | 0.0029 | U | 0.0031 | U | 0.003 | U |
| | - | - | | | | | | | | | | | | |
| Benzo(a)anthracene | 1 | 5.6 | 11 | 0.007 | U | 0.0066 U | 0.007 | U | 0.0065 | U | 0.007 | U | 0.0067 | U |
| Benzo(a)pyrene | 1 | 1 | 1.1 | 0.0077 | U | 0.0072 U | 0.0077 | U | 0.0071 | U | 0.0077 | U | 0.0074 | U |
| Benzo(b)fluoranthene | 1 | 5.6 | 11 | 0.0074 | U | 0.007 U | 0.0074 | U | 0.0069 | U | 0.0074 | U | 0.0071 | U |
| Benzo(g,h,i)perylene | 100 | 500 | 1000 | 0.011 | U | 0.01 U | 0.011 | U | 0.01 | U | 0.011 | U | 0.01 | U |
| Benzo(k)fluoranthene | 1 | 56 | 110 | 0.008 | U | 0.0076 U | 0.0081 | U | 0.0075 | U | 0.0081 | U | 0.0077 | U |
| 4-Bromophenyl phenyl ether | - | - | - | 0.0082 | U | 0.0078 U | 0.0082 | U | 0.0076 | U | 0.0082 | U | 0.0079 | U |
| Butyl benzyl phthalate | 100 | | | 0.019 | U | 0.018 U | 0.02 | U | 0.018 | U | 0.02 | U | 0.019 | U |
| 1,1'-Biphenyl | | | | 0.0067 | U | 0.0063 U | 0.0067 | U | 0.0062 | U | 0.0067 | U | 0.0064 | U |
| Benzaldehyde | - | - | - | 0.009 | U | 0.0085 U | 0.009 | U | 0.0084 | U | 0.009 | U | 0.0087 | U |
| 2-Chloronaphthalene | - | - | - | 0.0052 | U | 0.0049 U | 0.0052 | U | 0.0048 | U | 0.0052 | U | 0.005 | U |
| 4-Chloroaniline | 100 | | | 0.0096 | U | 0.009 U | 0.0096 | U | 0.0089 | U | 0.0096 | U | 0.0092 | U |
| Carbazole | - | - | - | 0.004 | U | 0.0038 U | 0.004 | U | 0.0037 | U | 0.004 | U | 0.0038 | U |
| Caprolactam | - | - | - | 0.023 | U | 0.022 U | 0.023 | U | 0.022 | U | 0.023 | U | 0.022 | U |
| Chrysene | 1 | 56 | 110 | 0.0058 | U | 0.0055 U | 0.0058 | U | 0.0054 | U | 0.0058 | U | 0.0056 | U |
| bis(2-Chloroethoxy)methane | - | - | - | 0.0082 | U | 0.0077 U | 0.0082 | U | 0.0076 | U | 0.0082 | U | 0.0079 | U |
| bis(2-Chloroethyl)ether | - | - | - | 0.015 | U | 0.014 U | 0.015 | U | 0.014 | U | 0.015 | U | 0.014 | U |
| bis(2-Chloroisopropyl)ether | - | - | - | 0.0083 | U | 0.0078 U | 0.0083 | U | 0.0077 | U | 0.0083 | U | 0.0079 | U |
| 4-Chlorophenyl phenyl ether | - | - | - | 0.0068 | U | 0.0064 U | 0.0068 | U | 0.0063 | U | 0.0068 | U | 0.0065 | U |
| 2,4-Dinitrotoluene | - | - | - | 0.0068 | U | 0.0064 U | 0.0068 | U | 0.0063 | U | 0.0068 | U | 0.0065 | U |
| 2,6-Dinitrotoluene | 1.03 | | | 0.0093 | U | 0.0088 U | 0.0093 | U | 0.0086 | U | 0.0093 | U | 0.0089 | U |
| 3,3'-Dichlorobenzidine | - | - | - | 0.024 | U | 0.022 U | 0.024 | U | 0.022 | U | 0.024 | U | 0.023 | U |
| 1,4-Dioxane | 9.8 | 130 | 250 | 0.024 | U | 0.023 U | 0.024 | U | 0.023 | U | 0.024 | U | 0.023 | U |
| Dibenzo(a,h)anthracene | 0.33 | 0.56 | 1.1 | 0.013 | U | 0.012 U | 0.013 | U | 0.012 | U | 0.013 | U | 0.012 | U |
| Dibenzofuran | 14 | 350 | 1000 | 0.005 | U | 0.0047 U | 0.005 | U | 0.0047 | U | 0.005 | U | 0.0048 | U |
| Di-n-butyl phthalate | 100 | 000 | 1000 | 0.0043 | U | 0.004 U | 0.0043 | U | 0.004 | U | 0.0043 | U | 0.0041 | U |
| Di-n-octyl phthalate | 100 | | | 0.0049 | U | 0.0046 U | 0.0049 | U | 0.0045 | U | 0.0049 | U | 0.0047 | U |
| Di-n-octyr phinalate Diethyl phthalate | 100 | | | 0.0049 | U | 0.0046 U 0.0043 U | 0.0049 | U | 0.0045 | U | 0.0049 | U | 0.0047 | U |
| | 100 | | | | U | | | | | U | | U | | U U |
| Dimethyl phthalate | 50 | | | 0.0052 | U | | 0.0052 | U | 0.0048 | | 0.0052 | U | 0.005 | U U |
| bis(2-Ethylhexyl)phthalate | | 500 | 4000 | | | 0.012 U | 1.68 | | 0.012 | U | 0.013 | | 0.012 | |
| Fluoranthene | 100 | 500 | 1000 | 0.0044 | U | 0.0041 U | 0.0044 | U | 0.0041 | U | 0.0044 | U | 0.0042 | U |
| Fluorene | 100 | 500 | 1000 | 0.0043 | U | 0.004 U | 0.0043 | U | 0.004 | U | 0.0043 | U | 0.0041 | U |
| Hexachlorobenzene | 0.41 | 6 | 12 | 0.0071 | U | 0.0067 U | 0.0071 | U | 0.0066 | U | 0.0071 | U | 0.0068 | U |
| Hexachlorobutadiene | - | - | - | 0.0096 | U | 0.009 U | 0.0096 | U | 0.0089 | U | 0.0096 | U | 0.0092 | U |
| Hexachlorocyclopentadiene | | | | 0.057 | U | 0.054 U | 0.057 | U | 0.053 | U | 0.057 | U | 0.055 | U |
| Hexachloroethane | - | - | - | 0.012 | U | 0.011 U | 0.012 | U | 0.011 | U | 0.012 | U | 0.011 | U |
| Indeno(1,2,3-cd)pyrene | 0.5 | 5.6 | 11 | 0.019 | U | 0.018 U | 0.019 | U | 0.017 | U | 0.019 | U | 0.018 | U |
| Isophorone | 100 | | | 0.0067 | U | 0.0064 U | 0.0068 | U | 0.0063 | U | 0.0068 | U | 0.0065 | U |
| 2-Methylnaphthalene | 0.41 | | | 0.0067 | U | 0.0064 U | 0.0068 | U | 0.0063 | U | 0.0068 | U | 0.0065 | U |
| 2-Nitroaniline | | | | 0.0082 | U | 0.0077 U | 0.0082 | U | 0.0076 | U | 0.0082 | U | 0.0079 | U |
| 3-Nitroaniline | | | | 0.01 | U | 0.0097 U | 0.01 | U | 0.0095 | U | 0.01 | U | 0.0098 | U |
| 4-Nitroaniline | - | - | - | 0.012 | U | 0.011 U | 0.012 | U | 0.011 | U | 0.012 | U | 0.012 | U |
| Naphthalene | 100 | 500 | 1000 | 0.0058 | U | 0.0054 U | 0.0058 | U | 0.0054 | U | 0.0058 | U | 0.0055 | U |
| Nitrobenzene | 3.7 | 69 | 140 | 0.011 | U | 0.011 U | 0.011 | U | 0.011 | U | 0.011 | U | 0.011 | U |
| N-Nitroso-di-n-propylamine | - | - | - | 0.011 | U | 0.01 U | 0.011 | U | 0.0099 | U | 0.011 | U | 0.01 | U |
| N-Nitrosodiphenylamine | | | | 0.019 | U | 0.018 U | 0.019 | U | 0.018 | U | 0.019 | U | 0.018 | U |
| | | | | | | | | | | | | | | |
| | 100 | 500 | 1000 | 0.004 | U | 0.0038 U | 0.004 | U | 0.0037 | U | 0.004 | U | 0.0038 | U |
| Phenanthrene Pyrene | 100 100 | 500 500 | 1000 1000 | 0.004 | U U | 0.0038 U 0.0043 U | 0.004 | U U | 0.0037 | U U | 0.004 | U U | 0.0038 | U U |

NOTE:

All values recorded in units of mg/kg

U - Value reported under the detection limit

J - Approximated value

Cells highlighted in green and italics indicate an exceedance of the parameter's Protection of Groundwater SCO, but not the Residential or Commercial SCO.

Cells highlighted in yellow and bold indicate exceedance of the parameter's Protection of Groundwater SCO and the Residential SCO, but not the Commercial SCO.

Cells highlighted in light red and bold italics indicate an exceedance of the parameter's Protection of Groundwater SCO, Residential SCO, and Commercial SCO.

<u>Table 3-5</u> Soil Sampling Results Summary (Inorganics)

| Analyte | NY SCO - Residential w/CP-51 (10/10) (6 NYCRR 375-6 12/06) | NY SCO - Commercial w/CP-51 (10/10) (6 NYCRR 375-6 12/06) | NY SCO - Industrial w/CP-51 (10/10) (6 NYCRR 375-6 12/06) | ನರ್ರ-1 1 ft 10/29/201 | 15 | ୪୫-1 5 ft 10/29/20 | 15 | SB-2 1 ft 10/29/20 | | SB-2 5 ft 10/29/20 | | ട്ടം. 1 ft 10/29/2 | | ട്ടം-3 5 ft 10/29/20 | |
|-------------------------|--|---|---|-----------------------------|----|--------------------------|----|--------------------------|---|--------------------------|---|--------------------------|---|----------------------------|---|
| Metals Analysis (mg/kg) | | | | | | | | | | | | | | | |
| Aluminum | | | | 6060 | | 4560 | | 4950 | | 4200 | | 4880 | | 5150 | |
| Antimony | | | | 2.2 | U | 2 | U | 2.2 | U | 2 | U | 2.3 | U | 2.1 | U |
| Arsenic | 16 | 16 | 16 | 2.2 | | 2 | U | 2.2 | U | 2 | U | 2.3 | U | 2.1 | U |
| Barium | 350 | 400 | 10000 | 25.2 | | 23.7 | | 22 | U | 21.6 | | 23.5 | | 23.8 | |
| Beryllium | 14 | 590 | 2700 | 0.44 | | 0.29 | | 0.28 | | 0.31 | | 0.32 | | 0.32 | |
| Cadmium | 2.5 | 9.3 | 60 | 0.55 | U | 0.5 | U | 0.55 | U | 0.5 | U | 0.57 | U | 0.52 | U |
| Calcium | | | | 1080 | | 1020 | | 1460 | | 868 | | 1000 | | 1210 | |
| Chromium | - | - | - | 18.6 | | 12.5 | | 13.8 | | 13.9 | | 15.5 | | 15.7 | |
| Cobalt | 30 | | | 9.4 | | 5.3 | | 32.3 | | 5.8 | | 8.4 | | 6.7 | |
| Copper | 270 | 270 | 10000 | 10.2 | | 9.3 | | 8.1 | | 8.7 | | 9.5 | | 9.4 | |
| Iron | 2000 | | | 12700 | | 9740 | | 10400 | | 9930 | | 11100 | | 10600 | |
| Lead | 400 | 1000 | 3900 | 5.6 | | 4.2 | | 4.4 | | 3.8 | | 5 | | 4.9 | |
| Magnesium | - | - | - | 2800 | | 2450 | | 2780 | | 2090 | | 2120 | | 2490 | |
| Manganese | 2000 | 10000 | 10000 | 272 | | 244 | | 210 | | 239 | | 257 | | 265 | |
| Mercury | 0.81 | 2.8 | 5.7 | 0.035 | U | 0.032 | U | 0.035 | U | 0.033 | U | 0.036 | U | 0.031 | U |
| Nickel | 140 | 310 | 10000 | 48.6 | | 50.7 | | 37 | | 41.5 | | 43 | | 46.1 | |
| Potassium | - | - | - | 1300 | | 990 | U | 1100 | U | 1000 | U | 1100 | U | 1000 | U |
| Selenium | 36 | 1500 | 6800 | 2.2 | U | 2 | U | 2.2 | U | 2 | U | 2.3 | U | 2.1 | U |
| Silver | 36 | 1500 | 6800 | 0.64 | | 0.5 | U | 1 | | 0.5 | U | 0.57 | U | 0.52 | U |
| Sodium | - | - | - | 1100 | U | 990 | U | 1100 | U | 1000 | U | 1100 | U | 1000 | U |
| Thallium | | | | 1.1 | U | 0.99 | U | 1.1 | U | 1 | U | 1.1 | U | 1 | U |
| Vanadium | 100 | | | 20.7 | | 16.1 | | 15.2 | | 14 | | 15.9 | | 16.1 | |
| Zinc | 2200 | 10000 | 10000 | 24 | | 20.8 | | 19.5 | | 19.1 | | 20.8 | | 20.3 | |

All values recorded in units of mg/kg

U - Value reported under the detection limit

Cells highlighted in green and italics indicate an exceedance of the parameter's Industrial SCO, but not the Residential or Commercial SCO.

Cells highlighted in yellow and bold indicate exceedance of the parameter's Industrial and Commercial SCO, but not the Residential SCO.

Cells highlighted in light red and bold italics indicate an exceedance of the parameter's Protection of Residential SCO, Commercial SCO, and Industrial SCO.

<u>Table 3-6</u> Groundwater Sampling Results Summary (VOCs)

| Analyte | NYSDEC Class | | Febru | uary 2016 Inve | estigation Su | mmary | | March 20 | 016 Suppleme | ental Data | Jul | y 2016 Suppl | emental Dat | a |
|-------------------------------|------------------------------------|--------------------|--------------------|-------------------|-------------------|--------------------|--------------------|--------------------|-------------------|-------------------|-------------------|------------------|------------------|-------------------|
| VOCs | GA Groundwater Quality Standard | MW-1I 2/10/2016 | MW-1S 2/10/2016 | MW-2 2/10/2016 | MW-3 2/10/2016 | MW-4S 2/10/2016 | MW-4I 2/10/2016 | MW-5S 3/18/2016 | MW-6 3/18/2016 | MW-7 3/18/2016 | MW-5I 7/8/2016 | MW-8 7/8/2016 | MW-9 7/8/2016 | MW-10 7/8/2016 |
| | | 2/10/2010 | 2/10/2010 | | 2/10/2010 | | 2/10/2010 | | 5/10/2010 | 3/10/2010 | | | | |
| Acetone | - 5 | | 1.6 U | | | | | 16 U | 1.6 U | | 3.8 U | 3.8 U | 3.8 U | 3.8 U |
| Acrolein | 5 5 | 1.6 U 2.6 U | | | 1.6 U 2.6 U | | 1.6 U | | | | | | | |
| Acrylonitrile | 5 | 2.0 0 | 2.6 U 0.1 U | | | 2.0 0 | 2.6 U | 26 U 1 U | | | | | | |
| Benzene Bromochloromethane | 5 | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 0.1 U | 1 0 | 0.1 U | 0.1 U | 0.14 U 0.46 U | | 0.66 0.46 U | 0.14 U 0.46 U |
| Bromodichloromethane | J | 0.1 U | 0.1 U | 0.1 U | 0.1 U | | 0.1 U | 1 U | 0.1 U | | | | | 0.46 U 0.55 U |
| | - | 0.1 U | 0.1 U | | 0.1 U | .. | 0.1 U | 1.7 U | | 0.1 U 0.17 U | 0.55 U 0.34 U | | 0.55 U 0.34 U | 0.35 U 0.34 U |
| Bromoform Bromomethane | 5 | 0.17 U | 0.17 U | 0 0 | 0.17 U | 0.17 U | 0.17 U | 5.7 U | | 0.17 U | 0.34 U | | 0.34 U 0.46 U | 0.34 U 0.46 U |
| 2-Butanone (MEK) | J | 0.57 0 | 0.57 0 | 0.57 0 | 0.57 0 | | 0.57 0 | | 0.57 0 | 0.57 0 | 1.9 U | 1.9 U | 1.9 U | 1.9 U |
| Carbon disulfide | | | | | | | | | | | 0.33 U | | 0.33 U | 0.33 U |
| Carbon tetrachloride | 5 | 0.096 U | 0.096 U | 0.096 U | 0.096 U | 0.096 U | 0.096 U | 0.96 U | 0.096 U | 0.096 U | 0.53 U | 0.53 U | 0.53 U 0.54 U | 0.33 U 0.54 U |
| Chlorobenzene | 5 | 0.090 U | 0.090 U | 0.090 U | 0.090 U | 0.090 U | 0.090 U | 0.90 U | | 0.090 U | 0.34 0 0.17 U | 0.34 0 0.17 U | 0.34 U | 0.34 U |
| Chloroethane | 5 | 0.093 0 0.21 U | 0.093 0 0.21 U | | 0.093 U 0.21 U | 0.093 U 0.21 U | 0.093 U 0.21 U | 2.1 U | | 0.093 0 0.21 U | 0.17 U | 0.17 U | 0.17 U 0.44 U | 0.17 U 0.44 U |
| 2-Chloroethyl vinyl ether | 5 | 0.21 U | 0.5 U | | 0.21 U | 0.21 U | 0.21 U | 5 U | | 0.21 U | 0.44 0 | 0.44 0 | 0.44 0 | 0.44 0 |
| Chloroform | - 7 | 0.89 J | 0.49 J | 0.62 J | 0.51 J | 0.56 J | 0.52 J | 0.91 U | | 1.5 | 0.78 J | 0.23 U | 0.23 U | 0.33 J |
| Chloromethane | 5 | 0.03 J | 0.43 J | 0.11 U | 0.11 U | 0.11 U | 0.02 J | 1.1 U | 0.031 U | 0.11 U | 0.96 U | 0.20 U | 0.96 U | 0.96 U |
| Cvclohexane | 5 | | 0.11 0 | 0.11 0 | 0.11 0 | | 0.11 0 | 1.1 0 | 0.11 0 | 0.11 0 | 0.30 U | 0.73 U | 0.30 U | 0.73 U |
| 1,2-Dibromo-3-chloropropane | 0.04 | | | | | | | | | | 0.69 U | 0.73 U | 0.69 U | 0.69 U |
| Dibromochloromethane | - | 0.15 U | 0.15 U | 0.15 U | 0.15 U | 0.15 U | 0.15 U | 1.5 U | 0.15 U | 0.15 U | 0.23 U | 0.23 U | 0.23 U | 0.23 U |
| 1.2-Dibromoethane | 0.0006 | 0.10 0 | 0.10 0 | 0.10 0 | 0.10 0 | | 0.10 0 | 1.0 0 | 0.10 0 | 0.10 0 | 0.22 U | 0.20 U | 0.20 U | 0.22 U |
| 1.2-Dichlorobenzene | 3 | 0.19 U | 0.72 J | 0.19 U | 0.19 U | 0.19 U | 0.19 U | 1.9 U | 0.19 U | 0.19 U | 0.22 0 | 0.22 U | 0.22 U | 0.23 U |
| 1.3-Dichlorobenzene | 3 | 0.19 U | 0.19 U | 0.19 U | 0.19 U | 0.19 U | 0.19 U | 1.9 U | | 0.19 U | 0.19 U | 0.19 U | 0.19 U | 0.19 U |
| 1.4-Dichlorobenzene | 3 | 0.11 U | 0.11 U | 0.11 U | 0.11 U | 0.11 U | 0.11 U | 1.1 U | | 0.11 U | 0.21 U | 0.21 U | 0.21 U | 0.21 U |
| Dichlorodifluoromethane | 5 | 0.29 U | 0.29 U | | 0.29 U | 0.29 U | 0.29 U | 2.9 U | 0.29 U | 0.29 U | 0.7 U | 0.7 U | 0.7 U | 0.7 U |
| 1,1-Dichloroethane | 5 | 0.12 U | 0.12 U | | 0.12 U | 0.12 U | 0.12 U | 1.2 U | | 0.12 U | 0.21 U | | 0.21 U | 0.21 U |
| 1,2-Dichloroethane | 0.6 | 0.09 U | 0.09 U | | 0.09 U | 0.09 U | 0.09 U | 0.9 U | | 0.09 U | 0.39 U | 0.39 U | 0.39 U | 0.39 U |
| 1,1-Dichloroethene | 5 | 0.16 U | 0.16 U | | 0.16 U | 0.16 U | 0.16 U | 1.6 U | 0.16 U | 0.16 U | 0.2 U | 0.2 U | 0.2 U | 0.2 U |
| cis-1,2-Dichloroethene | 5 | 0.12 U | 6.7 | 1.5 | 0.24 J | 0.12 U | 0.12 U | 687 | 2.1 | 0.34 J | 1.3 | 0.31 U | 1.1 | 2.9 |
| trans-1,2-Dichloroethene | 5 | 0.14 U | 0.14 U | | 0.14 U | 0.14 U | 0.14 U | 1.4 U | 0.14 U | 0.14 U | 0.36 U | 0.36 U | 0.36 U | 0.36 U |
| 1,2-Dichloropropane | 1 | 0.11 U | 0.11 U | | 0.11 U | 0.11 U | 0.11 U | 1.1 U | | 0.11 U | 0.33 U | | 0.33 U | 0.33 U |
| cis-1,3-Dichloropropene | _ | 0.12 U | 0.12 U | | 0.12 U | 0.12 U | 0.12 U | 1.2 U | | 0.12 U | 0.19 U | 0.19 U | 0.19 U | 0.19 U |
| trans-1,3-Dichloropropene | - | 0.15 U | 0.15 U | 0.15 U | 0.15 U | 0.15 U | 0.15 U | 1.5 U | | 0.15 U | 0.26 U | 0.26 U | 0.26 U | 0.26 U |
| Ethylbenzene | 5 | 0.22 U | | | 0.22 U | 0.22 U | 0.22 U | | | | | | 0.2 U | 0.2 U |
| Freon 113 | 5 | | | | | | | | | | 1.2 U | | 1.2 U | 1.2 U |
| 2-Hexanone | - | | | | | | | | | | 1.5 U | | 1.5 U | 1.5 U |
| Isopropylbenzene | 5 | | | | | | | | | | 0.16 U | | 0.16 U | 0.16 U |
| Methyl Acetate | - | | | | | | | | | | 1.5 U | | 1.5 U | 1.5 U |
| Methylcyclohexane | - | | | | | | | | | | 0.78 U | | 0.78 U | 0.78 U |
| Methyl Tert Butyl Ether | 10 | | | | | | | | | | 0.34 U | | 0.34 U | 0.34 U |
| 4-Methyl-2-pentanone(MIBK) | - | | | | | | | | | | 1.2 U | | 1.2 U | 1.2 U |
| Methylene chloride | 5 | 0.22 U | 0.22 U | 0.22 U | 0.22 U | 0.22 U | 0.22 U | 2.2 U | 0.22 U | 0.22 U | 0.35 U | | 0.35 U | 0.35 U |
| Styrene | 5 | | | | | | | | | | 0.27 U | | 0.27 U | 0.27 U |
| 1,1,2,2-Tetrachloroethane | 5 | 0.12 U | 0.12 U | 0.12 U | 0.12 U | 0.12 U | 0.12 U | 1.2 U | 0.12 U | 0.12 U | | | 0.39 U | 0.39 U |
| Tetrachloroethene | 5 | 0.72 J | 740 | 249 | 11.9 | 0.29 J | 0.26 J | 3490 | 3.7 | 4 | 18.7 | 0.52 J | 7.5 | 17.6 |



<u>Table 3-6</u> <u>Groundwater Sampling Results Summary (VOCs)</u>

| Analyte | NYSDEC Class | | Febru | uary 2016 Inve | estigation Su | mmary | | March 20 | 16 Suppleme | ntal Data | July | / 2016 Suppl | emental Dat | ta |
|------------------------|------------------------------------|--------------|--------------------|-------------------|-------------------|--------------------|--------------------|--------------------|-------------------|-------------------|-------------------|------------------|------------------|-------------------|
| VOCs | GA Groundwater Quality Standard | 141 4 4 - 11 | MW-1S 2/10/2016 | MW-2 2/10/2016 | MW-3 2/10/2016 | MW-4S 2/10/2016 | MW-4I 2/10/2016 | MW-5S 3/18/2016 | MW-6 3/18/2016 | MW-7 3/18/2016 | MW-5I 7/8/2016 | MW-8 7/8/2016 | MW-9 7/8/2016 | MW-10 7/8/2016 |
| Toluene | 5 | 0.25 U | 0.25 U | 0.25 U | 0.25 U | 0.25 U | 0.25 U | 2.5 U | 0.25 U | 0.25 U | 0.23 U | 0.23 U | 0.23 U | 0.23 U |
| 1,2,3-Trichlorobenzene | 5 | | | | | | | | | | 0.2 U | 0.2 U | 0.2 U | 0.2 U |
| 1,2,4-Trichlorobenzene | 5 | | | | | | | | | | 0.25 U | 0.25 U | 0.25 U | 0.25 U |
| 1,1,1-Trichloroethane | 5 | 0.086 U | 0.086 U | 0.086 U | 0.086 U | 0.086 U | 0.086 U | 0.86 U | 0.086 U | 0.086 U | 0.47 J | 0.22 U | 0.22 U | 0.22 U |
| 1,1,2-Trichloroethane | 1 | 0.13 U | 0.13 U | 0.13 U | 0.13 U | 0.13 U | 0.13 U | 1.3 U | 0.13 U | 0.13 U | 0.28 U | 0.28 U | 0.28 U | 0.28 U |
| Trichloroethene | 5 | 0.12 U | 9.7 | 2.4 | 0.3 J | 0.12 U | 0.12 U | 106 | 1.5 | 0.74 J | 2.1 | 0.26 U | 0.87 J | 3.4 |
| Trichlorofluoromethane | 5 | 0.2 U | 0.2 U | 0.2 U | 0.2 U | 0.2 U | 0.2 U | 2 U | 0.2 U | 0.2 U | 0.58 U | 0.58 U | 0.58 U | 0.58 U |
| Vinyl chloride | 2 | 0.13 U | 0.13 U | 0.13 U | 0.13 U | 0.13 U | 0.13 U | 1.3 U | 0.13 U | 0.13 U | 0.33 U | 0.33 U | 0.33 U | 0.33 U |
| Xylenes (total) | 5 | 0.22 U | 0.22 U | 0.22 U | 0.22 U | 0.22 U | 0.22 U | 2.2 U | 0.22 U | 0.22 U | 0.21 U | 0.21 U | 0.21 U | 0.21 U |

Notes:

All values recorded in units of ug/l

^a This compound outside control limits biased high in the associated BS.

U - Value reported under the detection limit

J - Approximated value

Cells highlighted in yellow / bold indicate the parameter exceeded the applicable NYSDEC Class GA Groundwater standard or guideline.



| Analyte GC/MS Semi-volatiles (EPA 625) | NYSDEC Class GA Groundwater Quality Standard | MW-2 2/10/2010 | 6 | MW-5 3/18/201 | 16 |
|---|--|-------------------|---|------------------|----|
| 2-Chlorophenol | - | 0.97 | U | 0.98 | U |
| 4-Chloro-3-methyl phenol | - | 1.2 | U | 1.2 | U |
| 2,4-Dichlorophenol | 1 | 1.6 | U | 1.6 | U |
| 2,4-Dimethylphenol | 1 | 1.7 | U | 1.7 | U |
| 2,4-Dinitrophenol | 1 | 0.91 | U | 0.91 | U |
| 4,6-Dinitro-o-cresol | | 0.74 | U | 0.74 | U |
| 2-Nitrophenol | - | 1.8 | U | 1.9 | U |
| 4-Nitrophenol | - | 0.86 | U | 0.86 | U |
| Pentachlorophenol | 1 | 1.9 | U | 1.9 | U |
| Phenol | 1 | 0.51 | U | 0.51 | U |
| 2,4,6-Trichlorophenol | - | 1.3 | U | 1.3 | U |
| Acenaphthene | - | 0.36 | U | 0.36 | U |
| Acenaphthylene | - | 0.39 | U | 0.30 | U |
| | | | - | | - |
| Anthracene | - | 0.41 | U | 0.41 | U |
| Benzidine | 5 | 0.29 | U | 0.29 | U |
| Benzo(a)anthracene | - | 0.36 | U | 0.37 | U |
| Benzo(a)pyrene | ND | 0.37 | U | 0.38 | U |
| Benzo(b)fluoranthene | - | 0.6 | U | 0.61 | U |
| Benzo(g,h,i)perylene | - | 0.43 | U | 0.43 | U |
| Benzo(k)fluoranthene | - | 0.43 | U | 0.43 | U |
| 4-Bromophenyl phenyl ether | - | 0.31 | U | 0.31 | U |
| Butyl benzyl phthalate | - | 0.61 | U | 0.61 | U |
| 2-Chloronaphthalene | - | 1 | U | 1 | U |
| 4-Chloroaniline | 5 | 0.41 | U | 0.41 | U |
| Chrysene | - | 0.26 | U | 0.26 | U |
| bis(2-Chloroethoxy)methane | 5 | 0.67 | U | 0.67 | U |
| bis(2-Chloroethyl)ether | 1 | 0.54 | C | 0.54 | U |
| bis(2-Chloroisopropyl)ether | 5 | 0.75 | U | 0.76 | U |
| 4-Chlorophenyl phenyl ether | - | 0.44 | U | 0.44 | U |
| 1,2-Dichlorobenzene | 3 | 0.21 | U | 0.21 | U |
| 1,2-Diphenylhydrazine | ND | 0.46 | U | 0.46 | U |
| 1,3-Dichlorobenzene | 3 | 0.16 | U | 0.16 | U |
| 1,4-Dichlorobenzene | 3 | 0.18 | U | 0.18 | U |
| 2,4-Dinitrotoluene | 5 | 0.88 | U | 0.88 | U |
| 2,6-Dinitrotoluene | 5 | 0.57 | U | 0.57 | U |
| 3,3'-Dichlorobenzidine | 5 | 1.2 | U | 1.3 | U |
| Dibenzo(a.h)anthracene | - | 0.56 | U | 0.56 | U |
| Di-n-butyl phthalate | 50 | 0.6 | U | 0.61 | U |
| Di-n-octyl phthalate | - | 0.58 | U | 0.58 | U |
| Diethyl phthalate | - | 0.4 | U | 0.4 | U |
| Dimethyl phthalate | - | 0.33 | U | 0.34 | U |
| bis(2-Ethylhexyl)phthalate | 5 | 0.67 | U | 0.68 | U |
| Fluoranthene | - | 0.87 | U | 0.08 | U |
| | - | 0.25 | U | | U |
| Huorene Hexachlorobenzene | 0.04 | 0.46 | U | 0.46 | U |
| | | | | | - |
| Hexachlorobutadiene | 0.5 | 0.18 | U | 0.18 | U |
| Hexachlorocyclopentadiene | 5 | 0.42 | U | 0.42 | U |
| Hexachloroethane | 5 | 0.29 | U | 0.29 | U |
| Indeno(1,2,3-cd)pyrene | - | 0.31 | U | 0.31 | U |
| Isophorone | - | 0.6 | U | 0.6 | U |
| Naphthalene | - | 0.33 | U | 0.33 | U |
| Nitrobenzene | 0.4 | 0.43 | U | 0.43 | U |
| n-Nitrosodimethylamine | - | 0.47 | U | 0.47 | U |
| N-Nitroso-di-n-propylamine | - | 0.48 | U | 0.48 | U |
| N-Nitrosodiphenylamine | - | 0.53 | U | 0.53 | U |
| Phenanthrene | - | 0.37 | U | 0.37 | U |
| Pyrene | - | 0.34 | U | 0.34 | U |
| 1,2,4-Trichlorobenzene | 5 | 0.35 | U | 0.35 | U |

<u>Table 3-7</u> <u>Groundwater Sampling Results Summary (SVOCs)</u>

Notes: All values recorded in units of ug/l ^a This compound outside control limits biased high in the associated BS.

U - Value reported under the detection limit

J - Approximated value

Cells highlighted in yellow / bold indicate the parameter exceeded the applicable NYSDEC Class GA Groundwater standard or guideline.

Table 3-8

Groundwater Sampling Results Summary (Pesticides / Herbicides)

| Analyte | NYSDEC Class GA Groundwater Quality Standard | MW-2 2/10/201 | 6 | MW-5 3/18/201 | 6 |
|---------------------------------------|--|------------------|---|------------------|---|
| GC Semi-volatiles (EPA 608) Aldrin | ND | 0.0017 | U | 0.0017 | U |
| alpha-BHC | 0.01 | 0.0017 | U | 0.0017 | U |
| beta-BHC | 0.04 | 0.0013 | U | 0.0013 | U |
| delta-BHC | 0.04 | 0.0016 | U | 0.0016 | U |
| gamma-BHC (Lindane) | 0.05 | 0.00086 | U | 0.00086 | U |
| Chlordane | 0.05 | 0.0000 | U | 0.0000 | U |
| Dieldrin | 0.004 | 0.00086 | U | 0.00086 | U |
| 4.4'-DDD | 0.3 | 0.00000 | U | 0.00000 | U |
| 4,4-DDE | 0.2 | 0.00084 | U | 0.00084 | U |
| 4,4'-DDT | 0.2 | 0.0025 | U | 0.0025 | U |
| Endrin | ND | 0.0025 | U | 0.0025 | U |
| Endosulfan sulfate | - | 0.0023 | U | 0.0023 | U |
| Endrin aldehyde | 5 | 0.0020 | U | 0.0032 a | U |
| Endosulfan-l | | 0.0011 | U | 0.0011 | U |
| Endosulfan-II | | 0.0016 | U | 0.0016 | U |
| Heptachlor | 0.04 | 0.0013 | U | 0.0013 | U |
| Heptachlor epoxide | 0.03 | 0.00074 | U | 0.00074 | U |
| Methoxychlor | 35 | 0.0034 | U | 0.0034 | U |
| Toxaphene | 0.06 | 0.047 | U | 0.047 | U |
| Aroclor 1016 | 0.09 | 0.047 | U | 0.048 | U |
| Aroclor 1221 | 0.09 | 0.24 | U | 0.24 | U |
| Aroclor 1232 | 0.09 | 0.2 | U | 0.2 | U |
| Aroclor 1242 | 0.09 | 0.083 | U | 0.085 | U |
| Aroclor 1248 | 0.09 | 0.077 | U | 0.079 | U |
| Aroclor 1254 | 0.09 | 0.055 | U | 0.056 | U |
| Aroclor 1260 | 0.09 | 0.059 | U | 0.06 | U |
| GC Semi-volatiles (SW846 8151) | | | | | |
| 2,4-D | 50 | 0.29 | U | 0.3 | U |
| 2,4,5-TP (Silvex) | 0.26 | 0.056 | U | 0.057 | U |
| 2,4,5-T | 35 | 0.057 | U | 0.058 | U |

Notes:

All values recorded in units of ug/l

^a This compound outside control limits biased high in the associated BS.

U - Value reported under the detection limit

J - Approximated value

Cells highlighted in yellow / bold indicate the parameter exceeded the applicable NYSDEC Class GA Groundwater standard or guideline.

<u>Table 3-9</u> <u>Groundwater Sampling Results Summary (Inorganics)</u>

| Analyte | NYSDEC Class GA Groundwater Quality Standard | MW-2 2/10/20 | | MW-3 3/18/20 | - |
|---------------------|--|-----------------|---|-----------------|---|
| Inorganics Analyses | | T | | | |
| Antimony | 3 | 6 | U | 6 | U |
| Arsenic | 25 | 3 | U | 3 | U |
| Beryllium | - | 1 | U | 1 | U |
| Cadmium | 5 | 3 | U | 3 | U |
| Chromium | 50 | 10 | U | 10 | U |
| Copper | 200 | 10 | U | 10 | U |
| Lead | 25 | 3 | U | 3 | U |
| Mercury | 0.7 | 0.2 | U | 0.2 | U |
| Nickel | 100 | 18.5 | | 10 | U |
| Selenium | 10 | 10 | U | 10 | U |
| Silver | 50 | 10 | U | 10 | U |
| Thallium | - | 2 | U | 2 | U |
| Zinc | - | 20 | U | 20 | U |

Notes:

All values recorded in units of ug/l

^a This compound outside control limits biased high in the associated BS.

U - Value reported under the detection limit

J - Approximated value

Cells highlighted in yellow / bold indicate the parameter exceeded the applicable NYSDEC Class GA Groundwater standard or guideline.

Figures

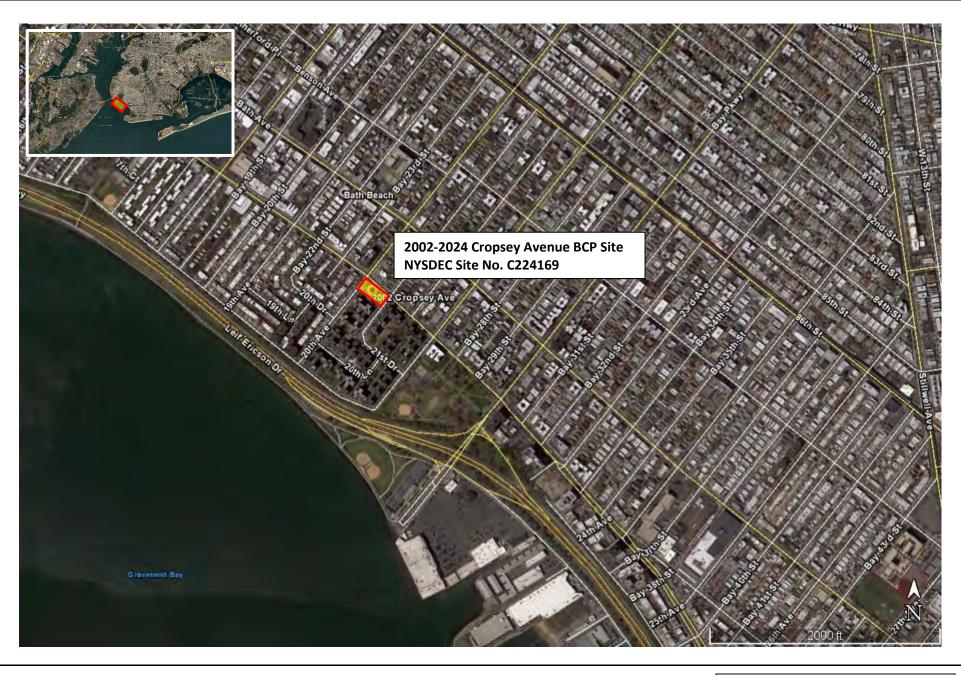




Figure 1 Site Location Map 2002-2024 Cropsey Avenue, Brooklyn, NY

 Client:
 20

 Project No.:
 85

 Project:
 20

 Date:
 Au









Figure 2 Remedial Investigation Areas Client: Project No.: Project: Date:

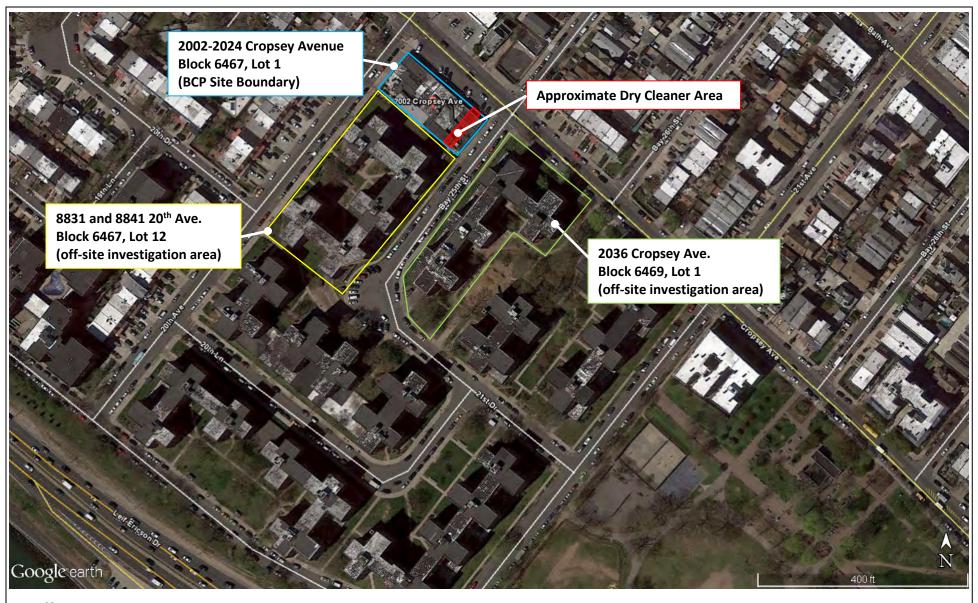






Figure 3 2002-2024 Cropsey Avenue Area Site Plan Client: Project No.: Project: Date:





Shallow Zone Monitoring Well (MW-#, MW-#S)

Intermediate Zone Monitoring Well (MW-#I)

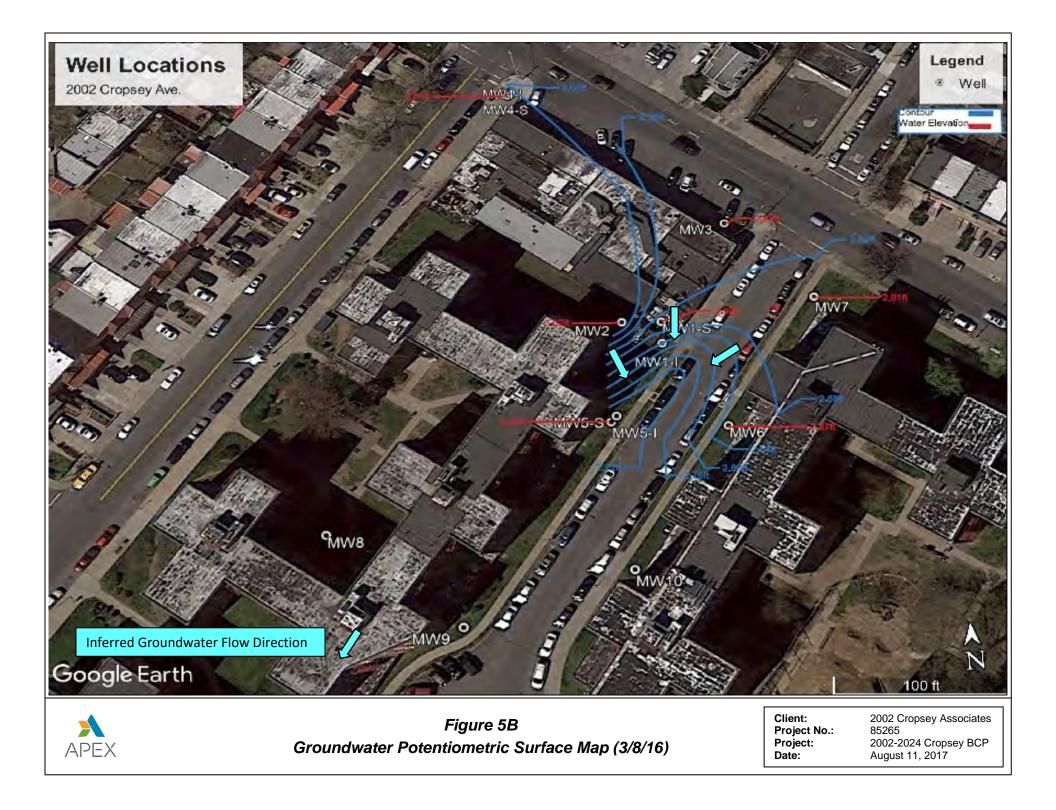


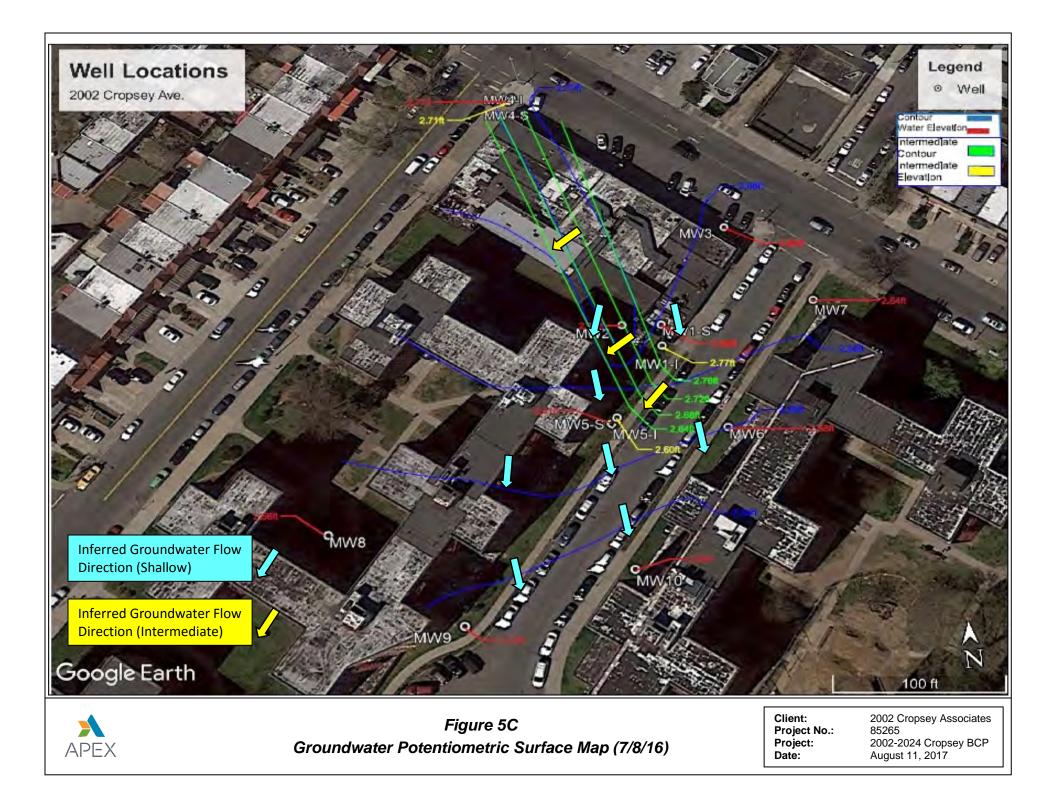
Figure 4 Monitoring Well Locations Client: Project No.: Project: Date:

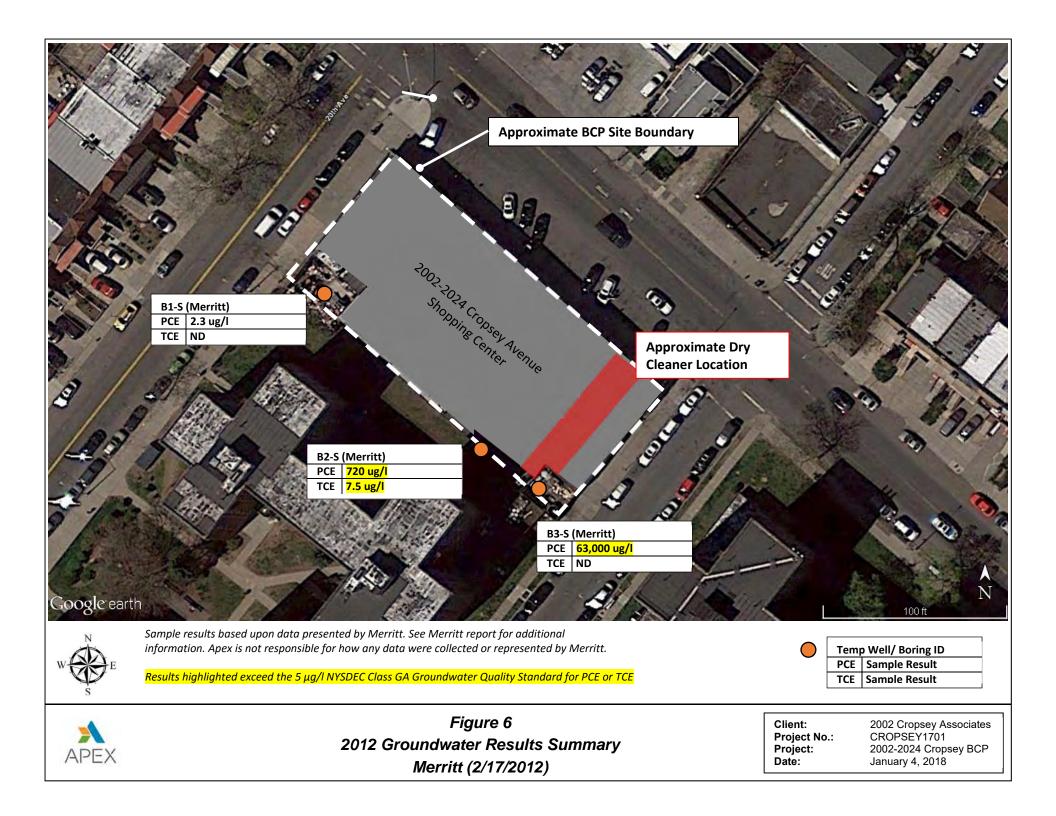


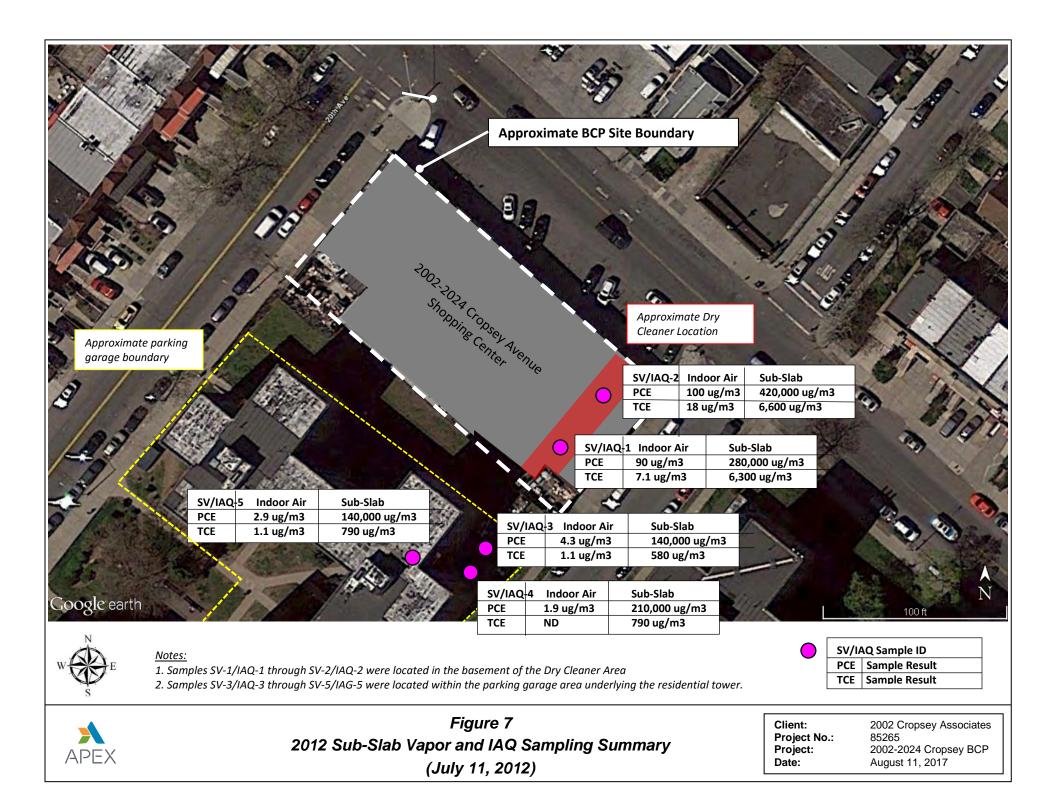


Figure 5A Groundwater Potentiometric Surface Map (7/27/12) Client: Project No.: Project: Date:









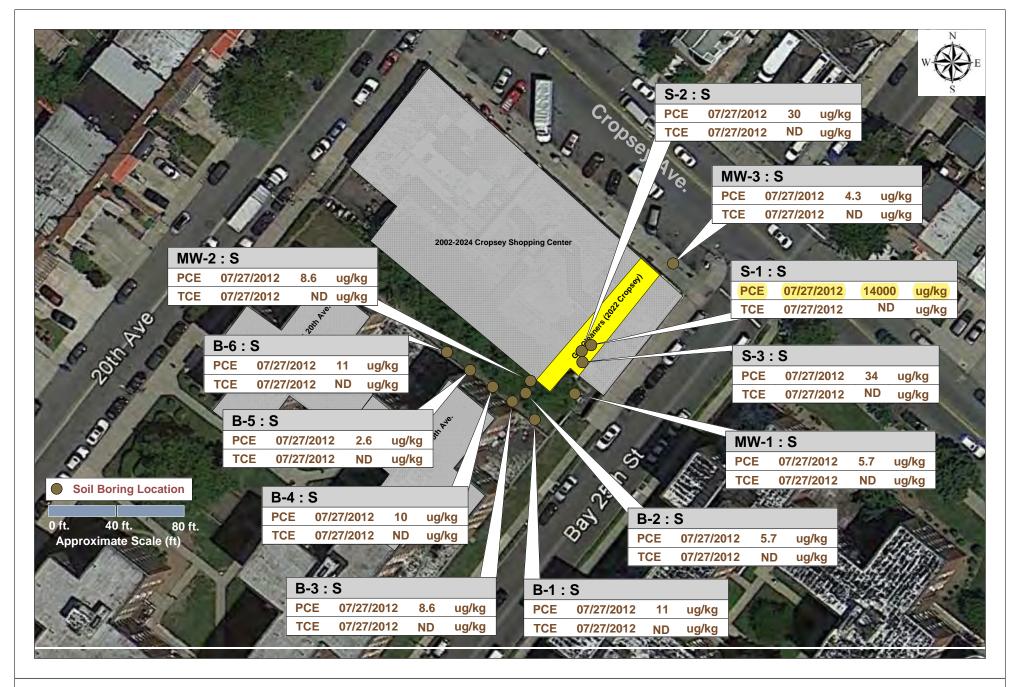
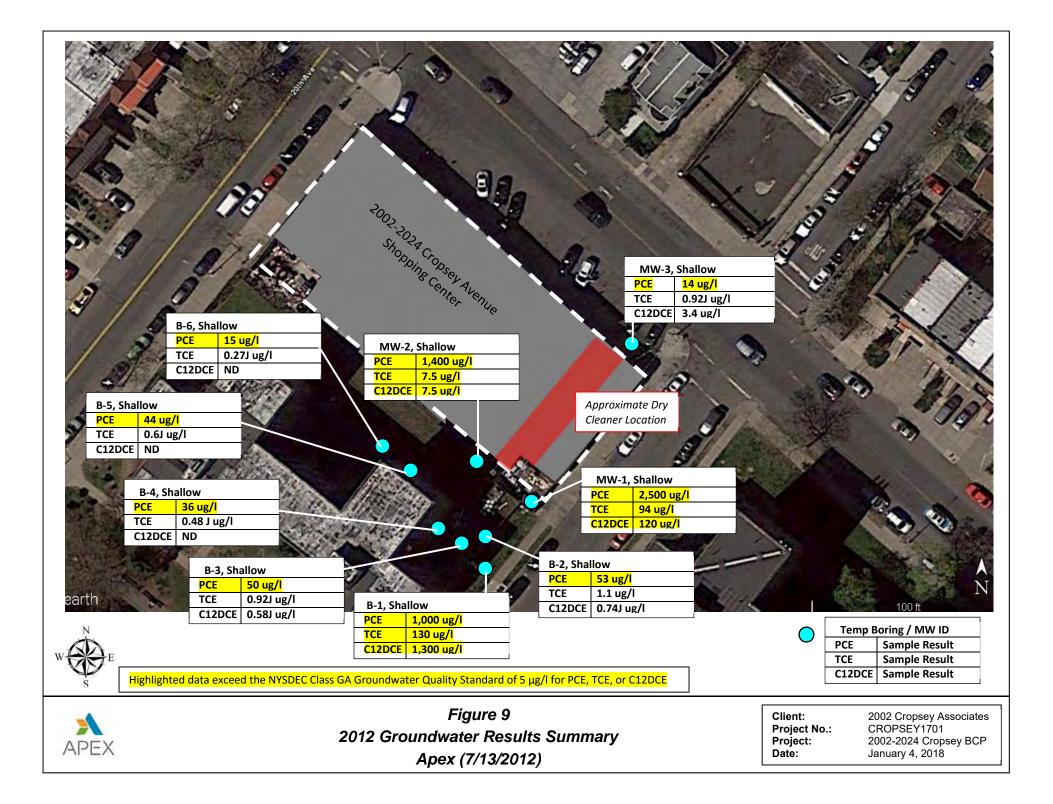




Figure 8 Soil Sample Results Summary (July 27, 2012)

Data highlighted exceeds NYSDEC Protection of GW SCO for that parameter



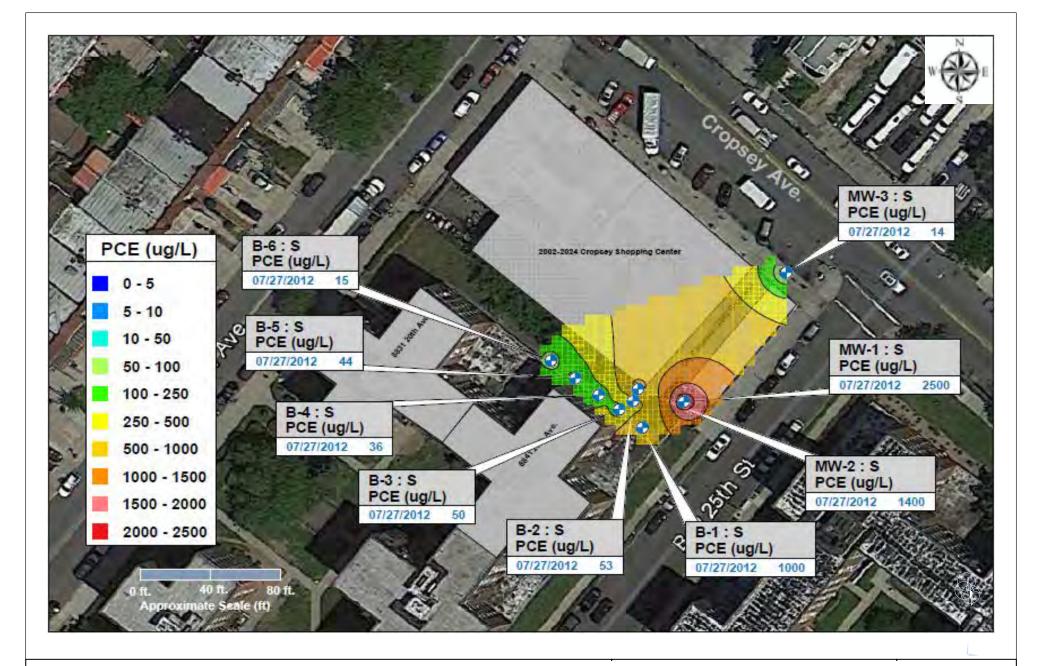




Figure 10 PCE Concentrations in Groundwater (July 27, 2012)







Figure 11 Membrane Interface Probe Investigation Locations Client: Project No.: Project: Date:

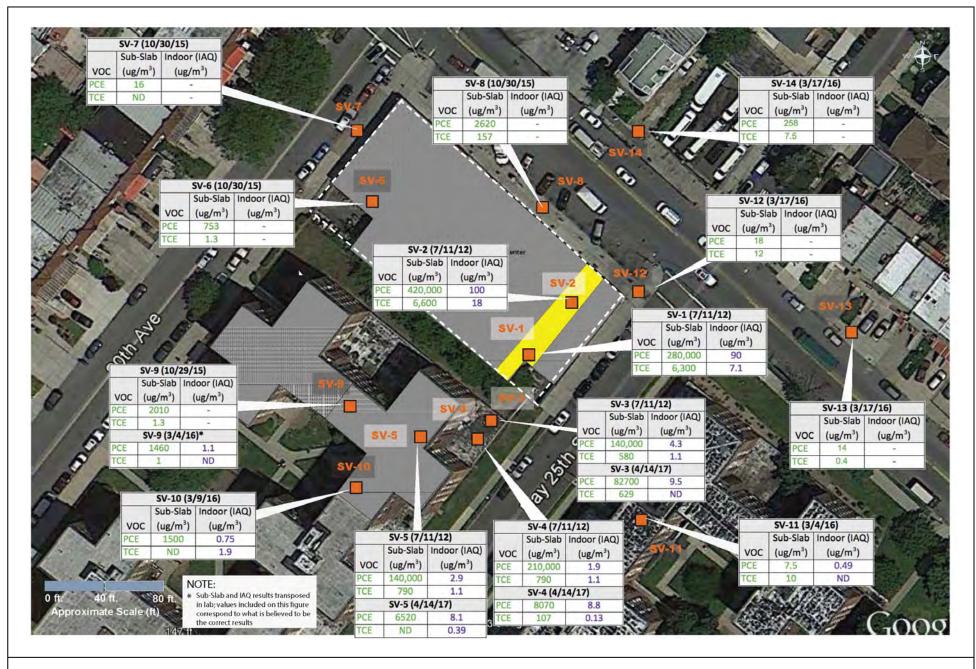
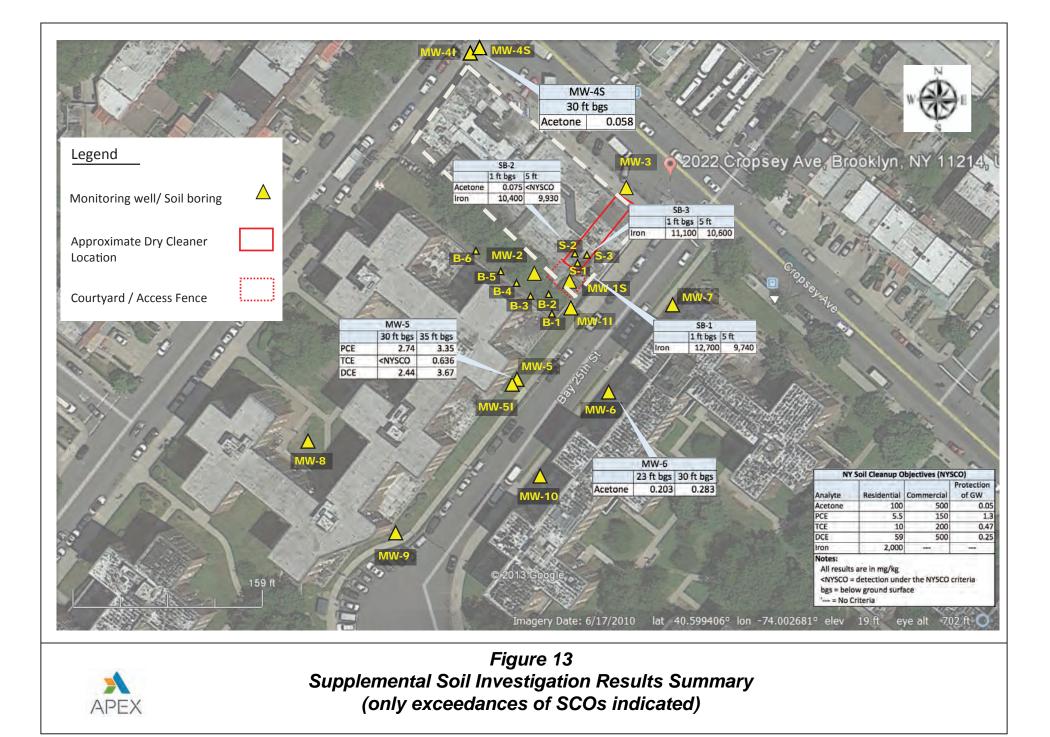




Figure 12 Soil Vapor / Indoor Air Sample Results Summary (All Events 2012 - 2017)



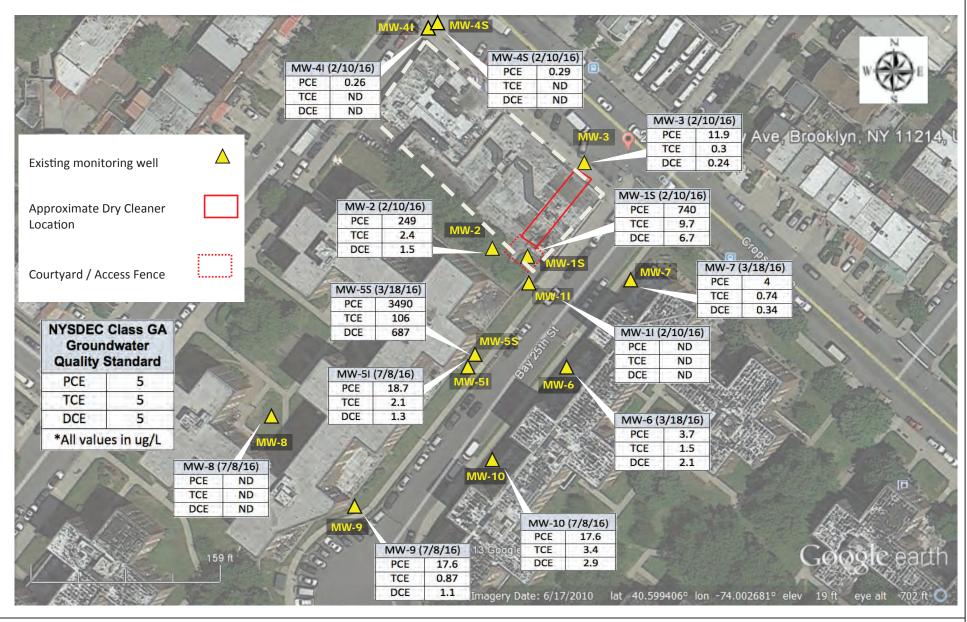
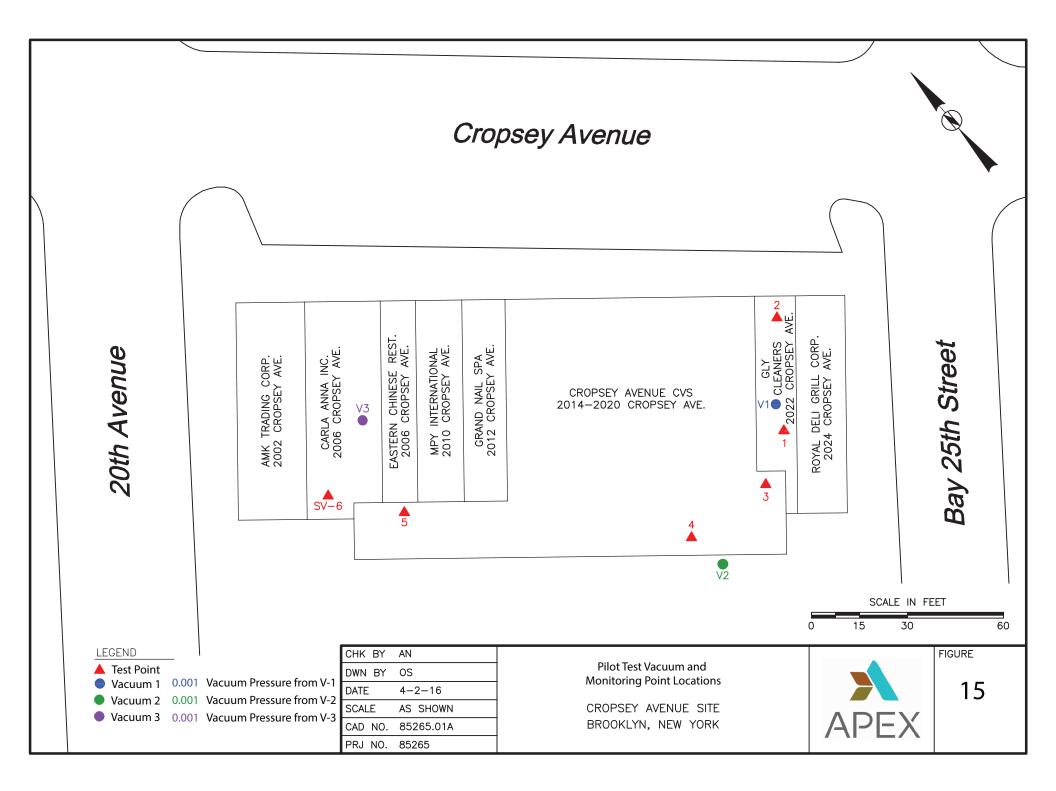
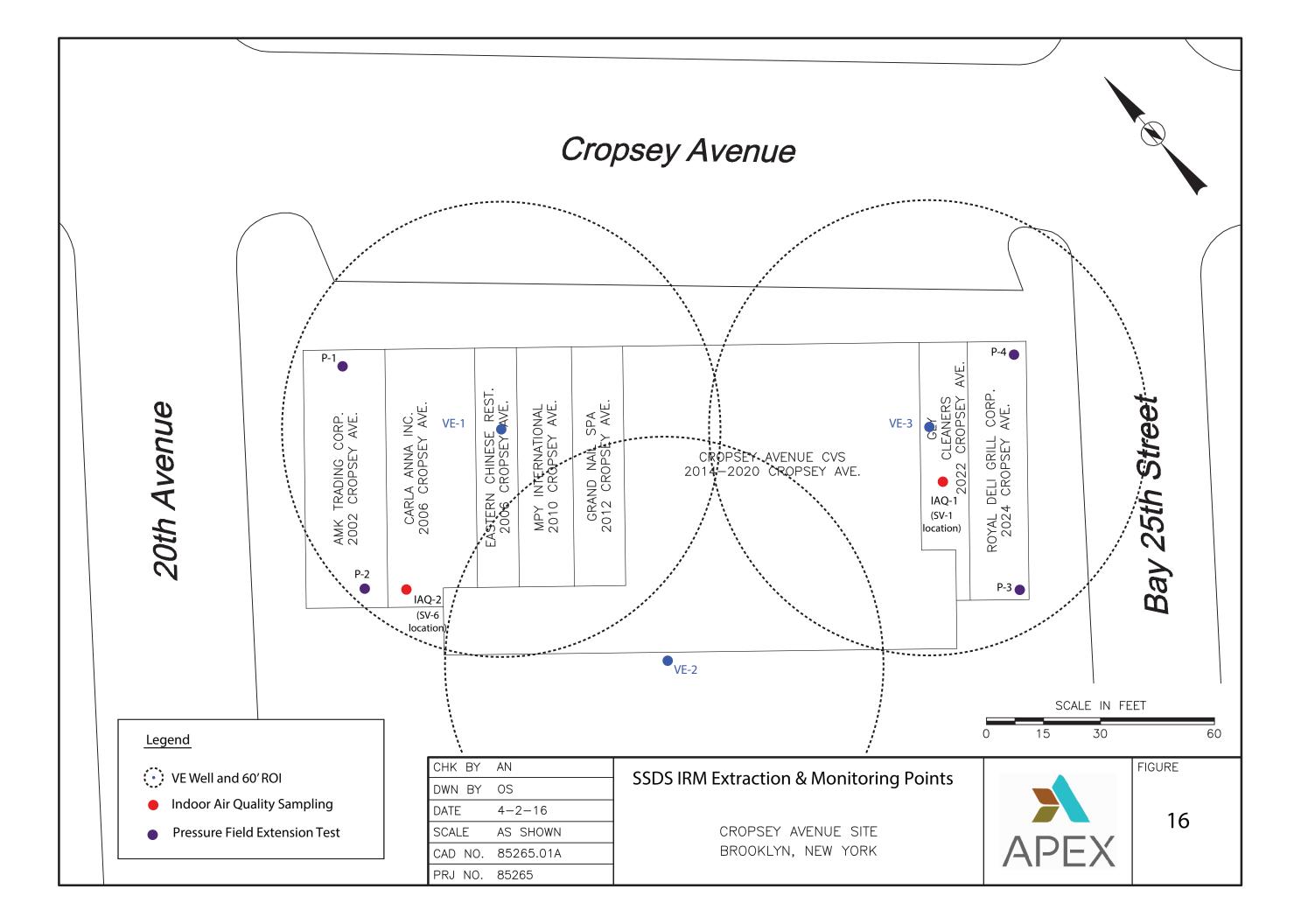


Figure 14 Groundwater Investigation Results Summary (only COCs indicated, 2016 composite data set)







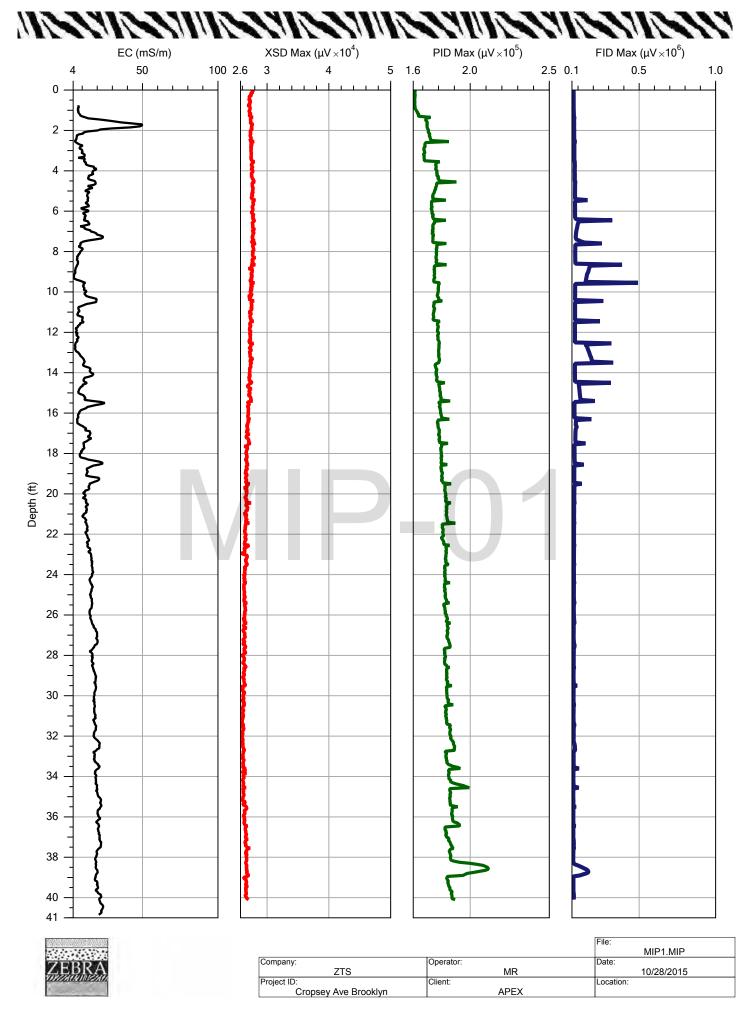
Appendix A

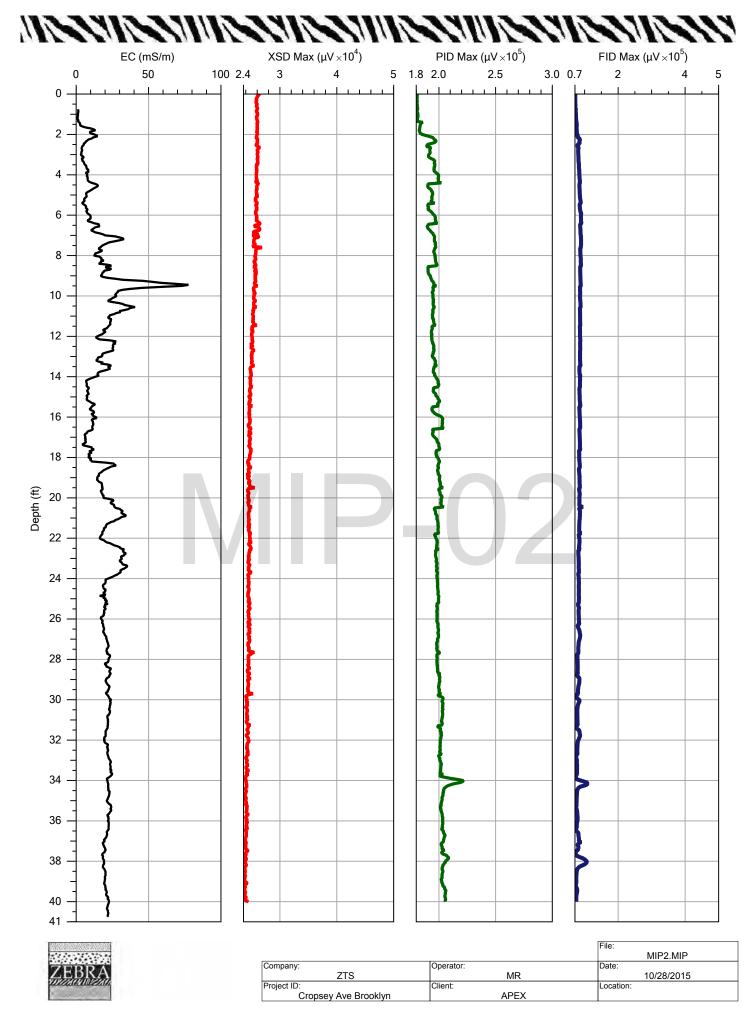
Apex Limited Phase II ESA Report

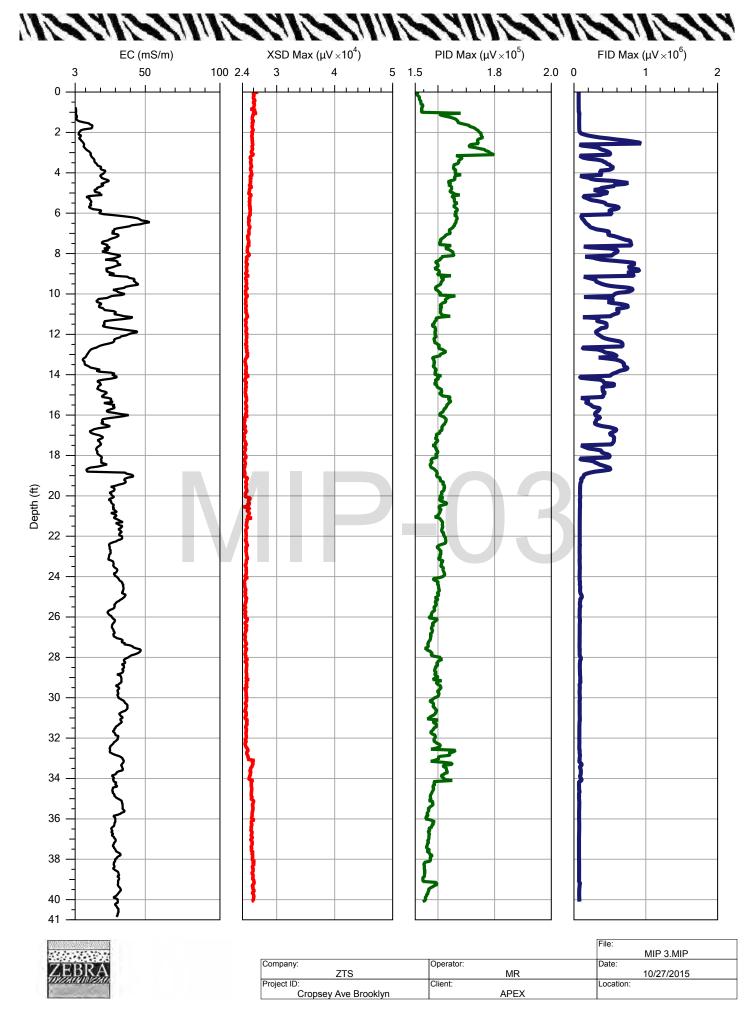
(incorporated by reference – under separate cover)

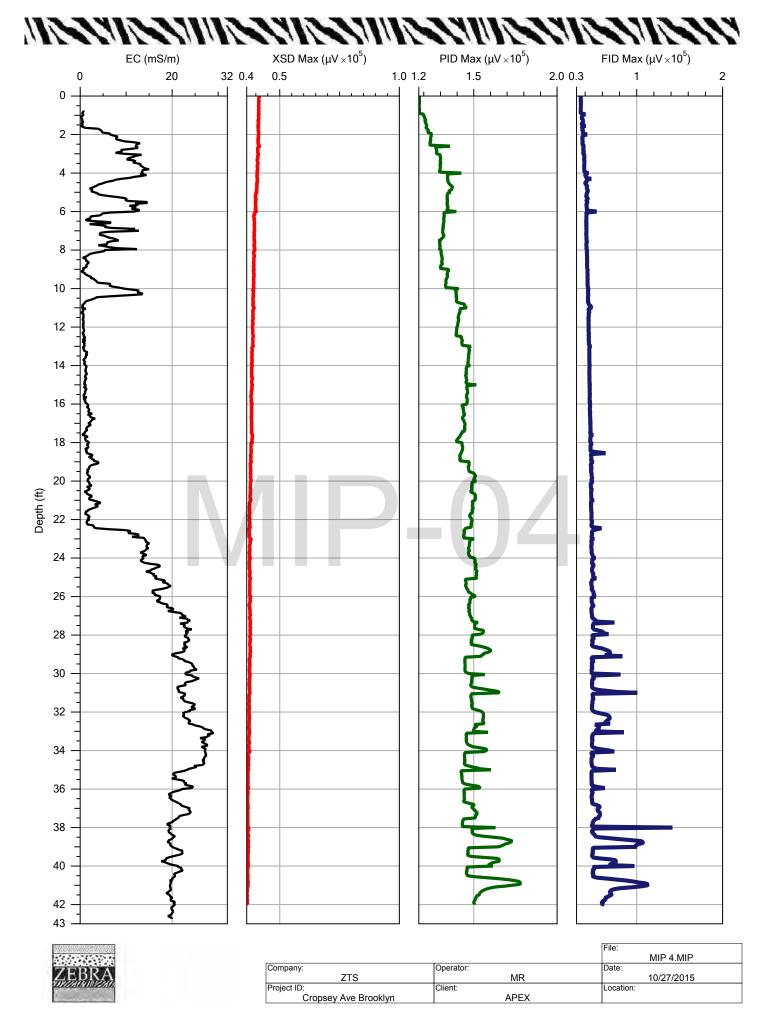
Appendix B

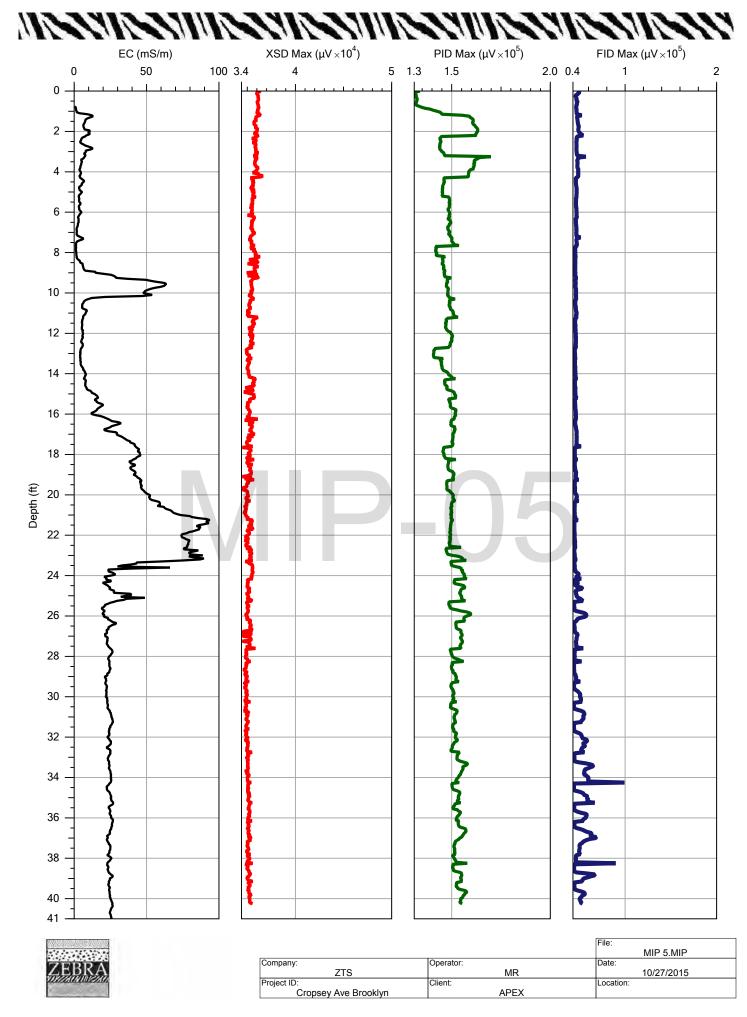
Membrane Interface Probe Investigation Data

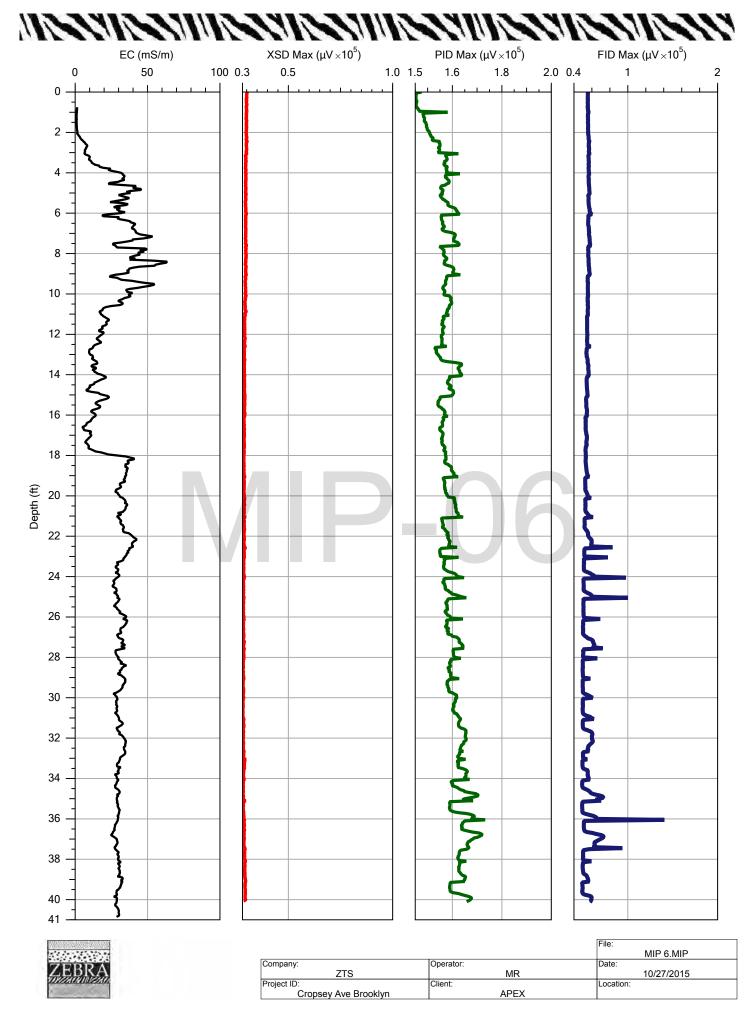




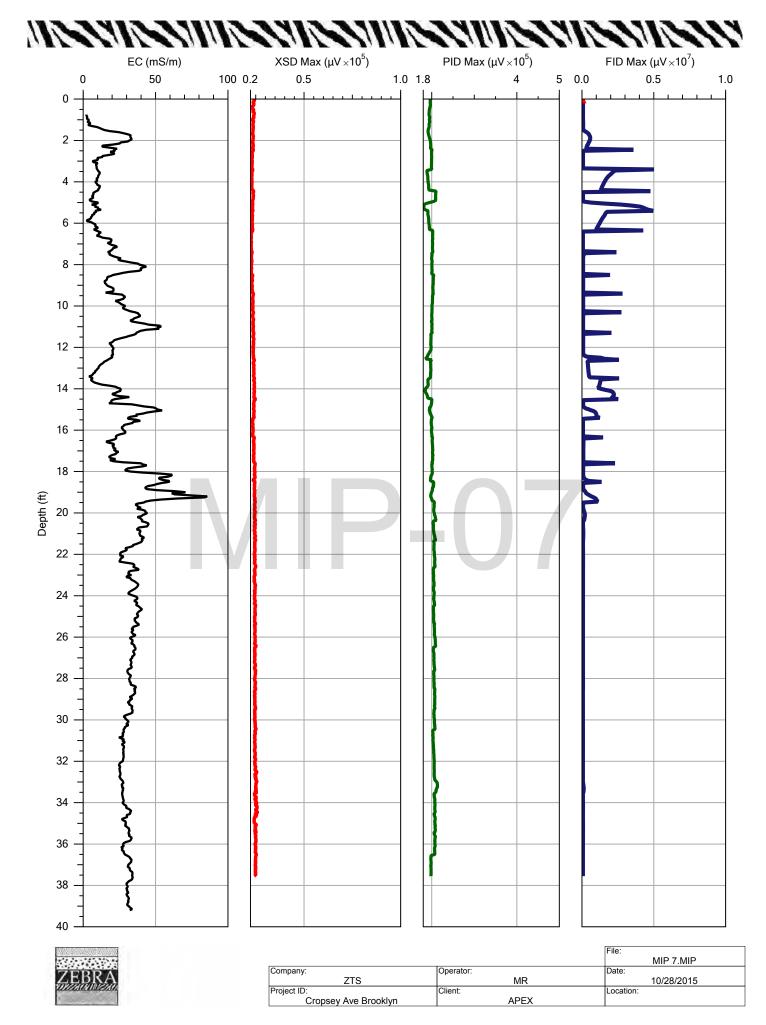




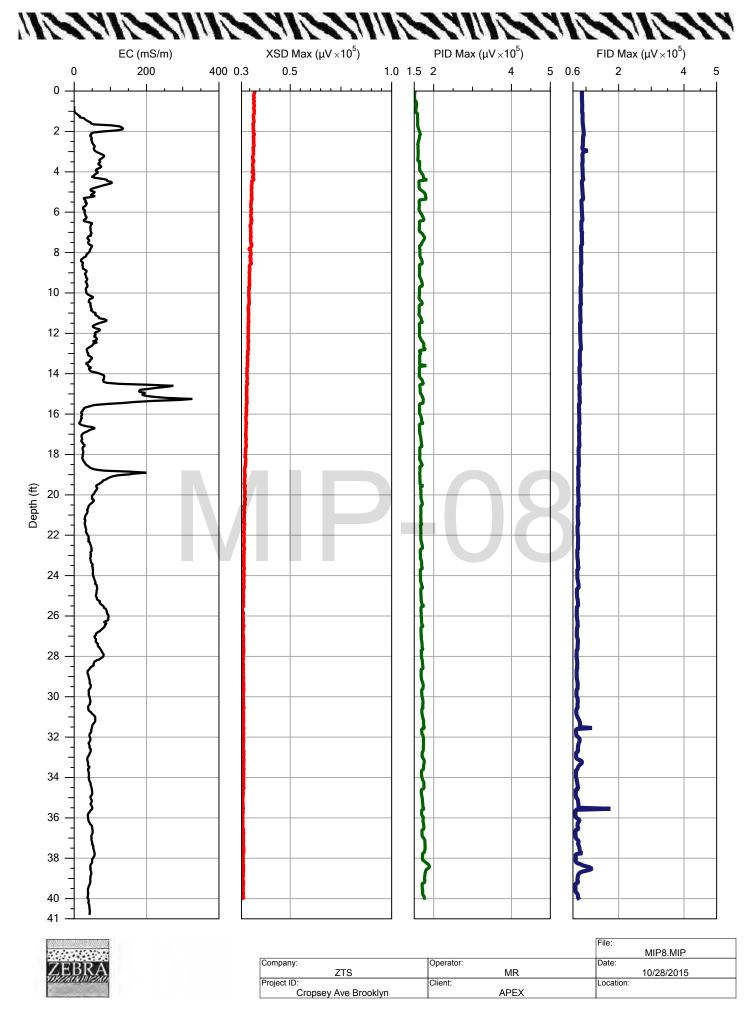




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Appendix C

Soil Vapor and Indoor Air Quality Data

Analytical Reports



Section 4

Sample Results

Report of Analysis





| Client San Lab Samj Matrix: Method: Project: | ple ID: JC751 AIR - TO-15 | 9-1 Soil Vapor | Comp. Summ ey Avenue, Bro | Date Sampled:10/29/15Date Received:10/31/15Percent Solids:n/a | | | |
|--|--|-------------------|------------------------------|---|-----------|------------|------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | 3W50810.D | 1 | 11/03/15 | YMH | n/a | n/a | V3W1930 |
| Run #2 | W52341.D | 1.48 | 11/04/15 | YMH | n/a | n/a | VW2090 |
| Run #1 Run #2 | Initial Volume 100 ml 40.0 ml | • | | | | | |

Report of Analysis

Page 1 of 3

4

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|---------|------------------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 280 ^a | 3.0 | 0.47 | ppbv | 665 ^a | 7.1 | 1.1 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.12 | ppbv | ND | 1.8 | 0.27 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 25.0 | 0.80 | 0.12 | ppbv | 79.9 | 2.6 | 0.38 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.13 | ppbv | ND | 5.4 | 0.87 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.082 | ppbv | ND | 8.3 | 0.85 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.087 | ppbv | ND | 3.1 | 0.34 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.081 | ppbv | ND | 3.5 | 0.35 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.10 | ppbv | ND | 4.1 | 0.52 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.80 | 0.11 | ppbv | ND | 2.5 | 0.34 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.13 | ppbv | ND | 3.7 | 0.60 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.086 | ppbv | ND | 2.1 | 0.23 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.80 | 0.12 | ppbv | ND | 3.9 | 0.59 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.12 | ppbv | ND | 1.7 | 0.25 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.13 | ppbv | ND | 4.1 | 0.67 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.098 | ppbv | ND | 5.0 | 0.62 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 1.2 | 0.80 | 0.13 | ppbv | 4.1 | 2.8 | 0.45 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.12 | ppbv | ND | 3.2 | 0.49 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | ND | 3.2 | 0.44 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.14 | ppbv | ND | 6.1 | 1.1 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.10 | ppbv | ND | 3.2 | 0.40 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.20 | ppbv | ND | 3.7 | 0.92 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.25 | ppbv | ND | 2.9 | 0.90 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.68 | 0.80 | 0.15 | ppbv J | 3.4 | 4.0 | 0.74 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.17 | ppbv | ND | 6.8 | 1.4 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.081 | ppbv | ND | 3.2 | 0.32 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.099 | ppbv | ND | 3.2 | 0.39 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.14 | ppbv | ND | 3.6 | 0.64 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | ND | 4.8 | 0.66 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.12 | ppbv | ND | 4.8 | 0.72 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.077 | ppbv | ND | 4.8 | 0.46 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.082 | ppbv | ND | 3.6 | 0.37 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



9 of 706 ACCUTEST JC7519

| Client Sample ID: | SV-09 | |
|-------------------|---|---------------|
| Lab Sample ID: | JC7519-1 Date Samp | led: 10/29/15 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A1090 Date Recei | ved: 10/31/15 |
| Method: | TO-15 Percent So | lids: n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|------------------|------|---------|---------|--------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 13.3 | 2.0 | 0.66 | ppbv | 25.1 | 3.8 | 1.2 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 23.0 | 0.80 | 0.19 | ppbv | 99.9 | 3.5 | 0.83 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 1.9 | 0.80 | 0.25 | ppbv | 6.8 | 2.9 | 0.90 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | 4.6 | 0.80 | 0.088 | ppbv | 23 | 3.9 | 0.43 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.11 | ppbv | ND | 6.1 | 0.84 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.10 | ppbv | ND | 5.6 | 0.70 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 13.2 | 0.80 | 0.12 | ppbv | 54.1 | 3.3 | 0.49 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.13 | ppbv | ND | 8.5 | 1.4 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 15.4 | 0.80 | 0.11 | ppbv | 54.3 | 2.8 | 0.39 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 0.80 | 0.18 | ppbv | ND | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 1.6 | 0.80 | 0.48 | ppbv | 3.9 | 2.0 | 1.2 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.54 | ppbv | ND | 2.8 | 1.9 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 13.1 | 0.80 | 0.19 | ppbv | 38.6 | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.80 | 0.11 | ppbv | ND | 3.3 | 0.45 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.10 | ppbv | ND | 2.9 | 0.36 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.12 | ppbv | ND | 3.3 | 0.49 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 2.0 | 0.32 | ppbv | ND | 3.4 | 0.55 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.10 | ppbv | ND | 3.4 | 0.43 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.13 | ppbv | ND | 4.4 | 0.71 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.12 | ppbv | ND | 5.5 | 0.82 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.14 | ppbv | ND | 4.4 | 0.76 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.18 | ppbv | ND | 5.9 | 1.3 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 7.4 | 0.80 | 0.092 | ppbv | 36 | 3.9 | 0.45 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 2.3 | 0.80 | 0.12 | ppbv | 11 | 3.9 | 0.59 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 1.2 | 0.80 | 0.085 | ppbv | 5.6 | 3.7 | 0.40 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 7.7 | 0.80 | 0.20 | ppbv | 23 | 2.4 | 0.61 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 297 ^a | 0.59 | 0.35 | ppbv | 2010 a | 4.0 | 2.4 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | 0.58 | 0.80 | 0.17 | ppbv J | 1.7 | 2.4 | 0.50 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 327 ^a | 3.0 | 0.30 | ppbv | 1230 a | 11 | 1.1 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 0.25 | 0.16 | 0.10 | ppbv | 1.3 | 0.86 | 0.54 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.081 | ppbv | ND | 4.5 | 0.46 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.13 | ppbv | ND | 2.0 | 0.33 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 82.0 | 0.80 | 0.17 | ppbv | 356 | 3.5 | 0.74 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 14.3 | 0.80 | 0.10 | ppbv | 62.1 | 3.5 | 0.43 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 96.4 | 0.80 | 0.10 | ppbv | 419 | 3.5 | 0.43 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | 2 | Limits | | | | | |
| 460-00-4 | 4-Brom | nofluorobenzene 105% | 96% | | 65-128% |) | | | | |

ND = Not detected MDL = Method Detection Limit

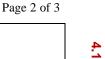
RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



10 of 706

ACCUTEST JC7519

| | Report of Analy | sis | Page 3 of 3 |
|---|--|-----------------------|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-09 JC7519-1 AIR - Soil Vapor Comp. Summa ID: A1090 TO-15 2002-2024 Cropsey Avenue, Brooklyn, NY | <u>*</u> 4 | |
| VOA TO15 List | | | |
| CAS No. MW | Compound Result RL | MDL Units Q Result RL | MDL Units |

(a) Result is from Run# 2

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



11 of 706 ACCUTEST JC7519

| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC751 AIR - TO-15 | 9-2 Soil Vapor | Comp. Summ ey Avenue, Bro | Date Sampled: 10/29/15 Date Received: 10/31/15 Percent Solids: n/a | | | |
|---|---|-------------------|------------------------------|--|-----------|------------|------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | 3W50811.D | 1 | 11/03/15 | YMH | n/a | n/a | V3W1930 |
| Run #2 | 3W50835.D | 1.48 | 11/05/15 | YMH | n/a | n/a | V3W1931 |
| Run #1 Run #2 | Initial Volum 100 ml 40.0 ml | 2 | | | | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|------|-------|-------|---|--------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 111 | 0.80 | 0.13 | ppbv | | 264 | 1.9 | 0.31 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.12 | ppbv | | ND | 1.8 | 0.27 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 22.4 | 0.80 | 0.12 | ppbv | | 71.6 | 2.6 | 0.38 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.13 | ppbv | | ND | 5.4 | 0.87 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.082 | ppbv | | ND | 8.3 | 0.85 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.087 | ppbv | | ND | 3.1 | 0.34 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.081 | ppbv | | ND | 3.5 | 0.35 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.10 | ppbv | | ND | 4.1 | 0.52 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.13 | ppbv | | ND | 3.7 | 0.60 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.086 | ppbv | | ND | 2.1 | 0.23 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | 0.56 | 0.80 | 0.12 | ppbv | J | 2.7 | 3.9 | 0.59 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.12 | ppbv | | ND | 1.7 | 0.25 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.13 | ppbv | | ND | 4.1 | 0.67 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.098 | ppbv | | ND | 5.0 | 0.62 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 3.2 | 0.80 | 0.13 | ppbv | | 11 | 2.8 | 0.45 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.12 | ppbv | | ND | 3.2 | 0.49 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | | ND | 3.2 | 0.44 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.14 | ppbv | | ND | 6.1 | 1.1 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.10 | ppbv | | ND | 3.2 | 0.40 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.20 | ppbv | | ND | 3.7 | 0.92 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.25 | ppbv | | ND | 2.9 | 0.90 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.73 | 0.80 | 0.15 | ppbv | J | 3.6 | 4.0 | 0.74 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.17 | ppbv | | ND | 6.8 | 1.4 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.081 | ppbv | | ND | 3.2 | 0.32 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.099 | ppbv | | ND | 3.2 | 0.39 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.14 | ppbv | | ND | 3.6 | 0.64 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | | ND | 4.8 | 0.66 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.12 | ppbv | | ND | 4.8 | 0.72 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.077 | ppbv | | ND | 4.8 | 0.46 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.082 | ppbv | | ND | 3.6 | 0.37 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

 $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



Page 1 of 3



E = Indicates value exceeds calibration range

| Report of Analysis | 5 |
|---------------------------|---|
|---------------------------|---|

| Client Sample ID: | SV-10 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7519-2 | Date Sampled: | 10/29/15 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A1045 | Date Received: | 10/31/15 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|------------------|------|---------|---------|-------------------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 8.5 | 2.0 | 0.66 | ppbv | 16 | 3.8 | 1.2 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 22.4 | 0.80 | 0.19 | ppbv | 97.3 | 3.5 | 0.83 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 1.1 | 0.80 | 0.25 | ppbv | 4.0 | 2.9 | 0.90 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | 5.1 | 0.80 | 0.088 | ppbv | 25 | 3.9 | 0.43 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.11 | ppbv | ND | 6.1 | 0.84 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.10 | ppbv | ND | 5.6 | 0.70 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 13.8 | 0.80 | 0.12 | ppbv | 56.6 | 3.3 | 0.49 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.13 | ppbv | ND | 8.5 | 1.4 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 19.3 | 0.80 | 0.11 | ppbv | 68.0 | 2.8 | 0.39 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 0.80 | 0.18 | ppbv | ND | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | ND | 0.80 | 0.48 | ppbv | ND | 2.0 | 1.2 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.54 | ppbv | ND | 2.8 | 1.9 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 8.1 | 0.80 | 0.19 | ppbv | 24 | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.80 | 0.11 | ppbv | ND | 3.3 | 0.45 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.10 | ppbv | ND | 2.9 | 0.36 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.12 | ppbv | ND | 3.3 | 0.49 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 2.0 | 0.32 | ppbv | ND | 3.4 | 0.55 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.10 | ppbv | ND | 3.4 | 0.43 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.13 | ppbv | ND | 4.4 | 0.71 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.12 | ppbv | ND | 5.5 | 0.82 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.14 | ppbv | ND | 4.4 | 0.76 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.18 | ppbv | ND | 5.9 | 1.3 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 8.2 | 0.80 | 0.092 | ppbv | 40 | 3.9 | 0.45 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 2.4 | 0.80 | 0.12 | ppbv | 12 | 3.9 | 0.59 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 2.8 | 0.80 | 0.085 | ppbv | 13 | 3.7 | 0.40 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 10.5 | 0.80 | 0.20 | ppbv | 31.8 | 2.4 | 0.61 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 370 ^a | 0.59 | 0.35 | ppbv | 2510 ^a | 4.0 | 2.4 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.80 | 0.17 | ppbv | ND | 2.4 | 0.50 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 198 a | 3.0 | 0.30 | ppbv | 746 ^a | 11 | 1.1 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 1.1 | 0.16 | 0.10 | ppbv | 5.9 | 0.86 | 0.54 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.081 | ppbv | ND | 4.5 | 0.46 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.13 | ppbv | ND | 2.0 | 0.33 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 81.5 | 0.80 | 0.17 | ppbv | 354 | 3.5 | 0.74 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 15.3 | 0.80 | 0.10 | ppbv | 66.5 | 3.5 | 0.43 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 96.9 | 0.80 | 0.10 | ppbv | 421 | 3.5 | 0.43 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | 2 | Limits | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 106% | 98% | | 65-128% | , | | | | |

ND = Not detectedMDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

4.2 4

Page 2 of 3

| | Report of Analy | vsis | Page 3 of 3 |
|---|--|-----------------------|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-10 JC7519-2 AIR - Soil Vapor Comp. Summa ID: A1045 TO-15 2002-2024 Cropsey Avenue, Brooklyn, NY | 4.2 4 | |
| VOA TO15 List CAS No. MW | Compound Result RL | MDL Units Q Result RL | MDL Units |

(a) Result is from Run# 2

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



14 of 706 ACCUTEST JC7519

| | | | | Repo | ort of A | nalysis | | Page 1 of 3 |
|---|-----------------------------|---|-------------------|--------------------------|------------------|-------------------------|--|------------------------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | | SV-11 JC7519-3 AIR - Soil Vapor Comp. Summa ID: A663 TO-15 2002-2024 Cropsey Avenue, Brooklyn, NY | | | | | Date Sampled: Date Received: Percent Solids: | |
| Run #1 ^a Run #2 | File ID W52368 | D | DF 1.38 | Analyzed 11/05/15 | Ву ҮМН | Prep Date n/a | Prep Batc n/a | h Analytical Batch VW2091 |
| Run #1 | Initial V 40.0 ml | olume | | | | | | |

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|-----|------|---------|--------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | ND | 2.8 | 0.44 | ppbv | ND | 6.7 | 1.0 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 2.8 | 0.43 | ppbv | ND | 6.2 | 0.95 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 12.6 | 2.8 | 0.41 | ppbv | 40.3 | 8.9 | 1.3 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 2.8 | 0.44 | ppbv | ND | 19 | 2.9 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 2.8 | 0.28 | ppbv | ND | 29 | 2.9 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 2.8 | 0.30 | ppbv | ND | 11 | 1.2 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 2.8 | 0.28 | ppbv | ND | 12 | 1.2 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 2.8 | 0.36 | ppbv | ND | 14 | 1.9 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 2.8 | 0.39 | ppbv | ND | 8.7 | 1.2 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 2.8 | 0.45 | ppbv | ND | 13 | 2.1 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 2.8 | 0.30 | ppbv | ND | 7.4 | 0.79 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 2.8 | 0.43 | ppbv | ND | 14 | 2.1 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 2.8 | 0.40 | ppbv | ND | 5.8 | 0.83 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 2.8 | 0.38 | ppbv | ND | 8.8 | 1.2 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 2.8 | 0.45 | ppbv | ND | 14 | 2.3 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 2.8 | 0.34 | ppbv | ND | 18 | 2.1 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 39.7 | 2.8 | 0.44 | ppbv | 137 | 9.6 | 1.5 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 2.8 | 0.43 | ppbv | ND | 11 | 1.7 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 2.8 | 0.39 | ppbv | ND | 11 | 1.5 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 2.8 | 0.49 | ppbv | ND | 22 | 3.8 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 2.8 | 0.36 | ppbv | ND | 11 | 1.5 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 2.8 | 0.69 | ppbv | ND | 13 | 3.2 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 2.8 | 0.87 | ppbv | ND | 10 | 3.1 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | ND | 2.8 | 0.51 | ppbv | ND | 14 | 2.5 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 2.8 | 0.57 | ppbv | ND | 24 | 4.9 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 2.8 | 0.28 | ppbv | ND | 11 | 1.1 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 2.8 | 0.34 | ppbv | ND | 11 | 1.3 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 2.8 | 0.49 | ppbv | ND | 13 | 2.2 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 2.8 | 0.38 | ppbv | ND | 17 | 2.3 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 2.8 | 0.41 | ppbv | ND | 17 | 2.5 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 2.8 | 0.26 | ppbv | ND | 17 | 1.6 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 2.8 | 0.28 | ppbv | ND | 13 | 1.3 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



SGS

| Client Sample ID: | SV-11 | | |
|-------------------|--|-------------|----------|
| Lab Sample ID: | JC7519-3 Date | Sampled: | 10/29/15 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A663 Date | Received: | 10/31/15 |
| Method: | TO-15 Perc | ent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|------|---------|-------|---|--------|-----|-----|-------|
| 64-17-5 | 46.07 | Ethanol | ND | 6.9 | 2.3 | ppbv | | ND | 13 | 4.3 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 0.92 | 2.8 | 0.66 | ppbv | J | 4.0 | 12 | 2.9 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | ND | 2.8 | 0.88 | ppbv | | ND | 10 | 3.2 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 2.8 | 0.30 | ppbv | | ND | 14 | 1.5 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 2.8 | 0.37 | ppbv | | ND | 21 | 2.8 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 2.8 | 0.35 | ppbv | | ND | 20 | 2.4 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 35.2 | 2.8 | 0.41 | ppbv | | 144 | 11 | 1.7 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 2.8 | 0.45 | ppbv | | ND | 30 | 4.8 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 61.2 | 2.8 | 0.38 | ppbv | | 216 | 9.9 | 1.3 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 2.8 | 0.61 | ppbv | | ND | 11 | 2.5 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | ND | 2.8 | 1.7 | ppbv | | ND | 6.9 | 4.2 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 2.8 | 1.8 | ppbv | | ND | 9.7 | 6.3 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | ND | 2.8 | 0.67 | ppbv | | ND | 8.3 | 2.0 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 2.8 | 0.37 | ppbv | | ND | 11 | 1.5 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 2.8 | 0.36 | ppbv | | ND | 10 | 1.3 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 2.8 | 0.42 | ppbv | | ND | 11 | 1.7 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 6.9 | 1.1 | ppbv | | ND | 12 | 1.9 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 2.8 | 0.35 | ppbv | | ND | 12 | 1.5 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 2.8 | 0.43 | ppbv | | ND | 15 | 2.3 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 2.8 | 0.42 | ppbv | | ND | 19 | 2.9 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 2.8 | 0.49 | ppbv | | ND | 15 | 2.7 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 2.8 | 0.61 | ppbv | | ND | 21 | 4.5 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | ND | 2.8 | 0.32 | ppbv | | ND | 14 | 1.6 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | 2.8 | 0.42 | ppbv | | ND | 14 | 2.1 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 12.8 | 2.8 | 0.29 | ppbv | | 59.8 | 13 | 1.4 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | ND | 2.8 | 0.69 | ppbv | | ND | 8.5 | 2.1 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | ND | 0.55 | 0.33 | ppbv | | ND | 3.7 | 2.2 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 2.8 | 0.60 | ppbv | | ND | 8.3 | 1.8 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 25.0 | 2.8 | 0.28 | ppbv | | 94.2 | 11 | 1.1 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 0.58 | 0.55 | 0.35 | ppbv | | 3.1 | 3.0 | 1.9 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 2.8 | 0.28 | ppbv | | ND | 16 | 1.6 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 2.8 | 0.45 | ppbv | | ND | 7.2 | 1.2 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 2.8 | 0.76 | ppbv | | ND | 9.8 | 2.7 | ug/m3 |
| | 106.2 | m,p-Xylene | 3.1 | 2.8 | 0.60 | ppbv | | 13 | 12 | 2.6 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 1.0 | 2.8 | 0.35 | ppbv | J | 4.3 | 12 | 1.5 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 4.2 | 2.8 | 0.35 | ppbv | | 18 | 12 | 1.5 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run | # 2 | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 96% | | | 65-128% | ó | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 3



16 of 706

| | Repo | ort of Analysis | 5 | | Page 3 of 3 |
|---|--|-----------------|--|--------|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-11 JC7519-3 AIR - Soil Vapor Comp. Summ TO-15 2002-2024 Cropsey Avenue, Bro | | Date Sampled: Date Received: Percent Solids: | | 4.3 |
| VOA TO15 List | | | | | |
| CAS No. MW | Compound | Result RL | MDL Units Q Resu | ılt RL | MDL Units |

(a) Diluted due to high concentration of non-target compound.

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



17 of 706 JC7519



Section 3 😡

Sample Results

Report of Analysis



| Client San Lab Samj Matrix: Method: Project: | ple ID: JC766 AIR - TO-15 | Soil Vapor | Comp. Summ ey Avenue, Bro | Date Sampled:10/30/15Date Received:11/03/15Percent Solids:n/a | | | | |
|--|--|------------|------------------------------|---|-----------|------------|--------------------|--|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batcl | h Analytical Batch | |
| Run #1 | 3W50844.D | 1 | 11/05/15 | YMH | n/a | n/a | V3W1931 | |
| Run #2 | 3W50862.D | 1.52 | 11/06/15 | YMH | n/a | n/a | V3W1932 | |
| Run #1 Run #2 | Initial Volume 100 ml 40.0 ml | | | | | | | |

Report of Analysis

Page 1 of 3

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VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|-------|---|------------------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 266 ^a | 3.0 | 0.48 | ppbv | | 632 ^a | 7.1 | 1.1 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.12 | ppbv | | ND | 1.8 | 0.27 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 1.9 | 0.80 | 0.12 | ppbv | | 6.1 | 2.6 | 0.38 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.13 | ppbv | | ND | 5.4 | 0.87 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.082 | ppbv | | ND | 8.3 | 0.85 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.087 | ppbv | | ND | 3.1 | 0.34 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.081 | ppbv | | ND | 3.5 | 0.35 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.10 | ppbv | | ND | 4.1 | 0.52 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.13 | ppbv | | ND | 3.7 | 0.60 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.086 | ppbv | | ND | 2.1 | 0.23 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | 2.5 | 0.80 | 0.12 | ppbv | | 12 | 3.9 | 0.59 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.12 | ppbv | | ND | 1.7 | 0.25 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.13 | ppbv | | ND | 4.1 | 0.67 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.098 | ppbv | | ND | 5.0 | 0.62 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 0.69 | 0.80 | 0.13 | ppbv | J | 2.4 | 2.8 | 0.45 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.12 | ppbv | | ND | 3.2 | 0.49 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | | ND | 3.2 | 0.44 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.14 | ppbv | | ND | 6.1 | 1.1 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.10 | ppbv | | ND | 3.2 | 0.40 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.20 | ppbv | | ND | 3.7 | 0.92 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.25 | ppbv | | ND | 2.9 | 0.90 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.60 | 0.80 | 0.15 | ppbv | J | 3.0 | 4.0 | 0.74 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.17 | ppbv | | ND | 6.8 | 1.4 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | 1.6 | 0.80 | 0.081 | ppbv | | 6.3 | 3.2 | 0.32 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | 1.6 | 0.80 | 0.099 | ppbv | | 6.3 | 3.2 | 0.39 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.14 | ppbv | | ND | 3.6 | 0.64 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | | ND | 4.8 | 0.66 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.12 | ppbv | | ND | 4.8 | 0.72 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.077 | ppbv | | ND | 4.8 | 0.46 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.082 | ppbv | | ND | 3.6 | 0.37 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



| Client Sample ID: | SV-8 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC7667-1 | Date Sampled: | 10/30/15 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A1046 | Date Received: | 11/03/15 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|------------------|------|---------|---------|--------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 8.5 | 2.0 | 0.66 | ppbv | 16 | 3.8 | 1.2 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 3.5 | 0.80 | 0.19 | ppbv | 15 | 3.5 | 0.83 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | ND | 0.80 | 0.25 | ppbv | ND | 2.9 | 0.90 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | 1.2 | 0.80 | 0.088 | ppbv | 5.9 | 3.9 | 0.43 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.11 | ppbv | ND | 6.1 | 0.84 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.10 | ppbv | ND | 5.6 | 0.70 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 3.7 | 0.80 | 0.12 | ppbv | 15 | 3.3 | 0.49 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.13 | ppbv | ND | 8.5 | 1.4 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 1.9 | 0.80 | 0.11 | ppbv | 6.7 | 2.8 | 0.39 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | 3.5 | 0.80 | 0.18 | ppbv | 14 | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 1.6 | 0.80 | 0.48 | ppbv | 3.9 | 2.0 | 1.2 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.54 | ppbv | ND | 2.8 | 1.9 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 12.5 | 0.80 | 0.19 | ppbv | 36.9 | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.80 | 0.11 | ppbv | ND | 3.3 | 0.45 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.10 | ppbv | ND | 2.9 | 0.36 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.12 | ppbv | ND | 3.3 | 0.49 | ug/m3 |
| 115-07-1 | 42 | Propylene | 12.9 | 2.0 | 0.32 | ppbv | 22.2 | 3.4 | 0.55 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.10 | ppbv | ND | 3.4 | 0.43 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.13 | ppbv | ND | 4.4 | 0.71 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.12 | ppbv | ND | 5.5 | 0.82 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.14 | ppbv | ND | 4.4 | 0.76 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.18 | ppbv | ND | 5.9 | 1.3 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 3.5 | 0.80 | 0.092 | ppbv | 17 | 3.9 | 0.45 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 1.2 | 0.80 | 0.12 | ppbv | 5.9 | 3.9 | 0.59 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 0.67 | 0.80 | 0.085 | ppbv J | 3.1 | 3.7 | 0.40 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 8.2 | 0.80 | 0.20 | ppbv | 25 | 2.4 | 0.61 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 387 ^a | 0.61 | 0.36 | ppbv | 2620 a | 4.1 | 2.4 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.80 | 0.17 | ppbv | ND | 2.4 | 0.50 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 15.5 | 0.80 | 0.081 | ppbv | 58.4 | 3.0 | 0.31 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 29.3 | 0.16 | 0.10 | ppbv | 157 | 0.86 | 0.54 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.081 | ppbv | ND | 4.5 | 0.46 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.13 | ppbv | ND | 2.0 | 0.33 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 13.3 | 0.80 | 0.17 | ppbv | 57.8 | 3.5 | 0.74 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 4.7 | 0.80 | 0.10 | ppbv | 20 | 3.5 | 0.43 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 17.9 | 0.80 | 0.10 | ppbv | 77.7 | 3.5 | 0.43 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | 2 | Limits | | | | | |
| 460-00-4 | 4-Brom | nofluorobenzene 105% | 96% | | 65-128% |) | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Page 2 of 3

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| | Report of Analysis | Page 3 of 3 | 3. <u>1</u> |
|---|--|-------------|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-8 JC7667-1 Date Sampled: 10/30/15 AIR - Soil Vapor Comp. Summa ID: A1046 Date Received: 11/03/15 TO-15 Percent Solids: n/a 2002-2024 Cropsey Avenue, Brooklyn, NY n/a | | ယ |
| VOA TO15 List | | , | |
| CAS No. MW | Compound Result RL MDL Units Q Result RL | MDL Un | nits |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



11 of 23 ACCUTEST JC7667

| Client Sa Lab Sam Matrix: Method: Project: | AIR - TO-15 | Soil Vapor | Comp. Summ ey Avenue, Bro | | 73 D P | Date Sampled:10/30/15Date Received:11/03/15Percent Solids:n/a | | | | | |
|--|----------------|------------|------------------------------|-----|-----------|---|------------------|--|--|--|--|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | | | | |
| Run #1 | 3W50845.D | 1 | 11/05/15 | YMH | n/a | n/a | V3W1931 | | | | |
| Run #2 | 3W50863.D | 1.52 | 11/06/15 | YMH | n/a | n/a | V3W1932 | | | | |
| | Initial Volume | • | | | | | | | | | |

Report of Analysis

Page 1 of 3

3.2

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VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|-------|---|--------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 515 ^a | 6.1 | 0.96 | ppbv | | 1220 a | 14 | 2.3 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.12 | ppbv | | ND | 1.8 | 0.27 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 1.7 | 0.80 | 0.12 | ppbv | | 5.4 | 2.6 | 0.38 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.13 | ppbv | | ND | 5.4 | 0.87 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.082 | ppbv | | ND | 8.3 | 0.85 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.087 | ppbv | | ND | 3.1 | 0.34 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.081 | ppbv | | ND | 3.5 | 0.35 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.10 | ppbv | | ND | 4.1 | 0.52 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.13 | ppbv | | ND | 3.7 | 0.60 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.086 | ppbv | | ND | 2.1 | 0.23 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.80 | 0.12 | ppbv | | ND | 3.9 | 0.59 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.12 | ppbv | | ND | 1.7 | 0.25 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.13 | ppbv | | ND | 4.1 | 0.67 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.098 | ppbv | | ND | 5.0 | 0.62 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 0.76 | 0.80 | 0.13 | ppbv | J | 2.6 | 2.8 | 0.45 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.12 | ppbv | | ND | 3.2 | 0.49 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | | ND | 3.2 | 0.44 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.14 | ppbv | | ND | 6.1 | 1.1 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.10 | ppbv | | ND | 3.2 | 0.40 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.20 | ppbv | | ND | 3.7 | 0.92 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.25 | ppbv | | ND | 2.9 | 0.90 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.52 | 0.80 | 0.15 | ppbv | J | 2.6 | 4.0 | 0.74 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.17 | ppbv | | ND | 6.8 | 1.4 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.081 | ppbv | | ND | 3.2 | 0.32 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.099 | ppbv | | ND | 3.2 | 0.39 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.14 | ppbv | | ND | 3.6 | 0.64 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | | ND | 4.8 | 0.66 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.12 | ppbv | | ND | 4.8 | 0.72 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.077 | ppbv | | ND | 4.8 | 0.46 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.082 | ppbv | | ND | 3.6 | 0.37 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



12 of 23

| Client Sample ID: | SV-7 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7667-2 | Date Sampled: | 10/30/15 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A873 | Date Received: | 11/03/15 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|------|---------|---------|--------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 8.6 | 2.0 | 0.66 | ppbv | 16 | 3.8 | 1.2 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 4.3 | 0.80 | 0.19 | ppbv | 19 | 3.5 | 0.83 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 2.8 | 0.80 | 0.25 | ppbv | 10 | 2.9 | 0.90 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | 1.4 | 0.80 | 0.088 | ppbv | 6.9 | 3.9 | 0.43 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.11 | ppbv | ND | 6.1 | 0.84 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.10 | ppbv | ND | 5.6 | 0.70 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 4.7 | 0.80 | 0.12 | ppbv | 19 | 3.3 | 0.49 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.13 | ppbv | ND | 8.5 | 1.4 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 2.1 | 0.80 | 0.11 | ppbv | 7.4 | 2.8 | 0.39 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | 5.0 | 0.80 | 0.18 | ppbv | 20 | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | ND | 0.80 | 0.48 | ppbv | ND | 2.0 | 1.2 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.54 | ppbv | ND | 2.8 | 1.9 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 21.7 | 0.80 | 0.19 | ppbv | 64.0 | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.80 | 0.11 | ppbv | ND | 3.3 | 0.45 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.10 | ppbv | ND | 2.9 | 0.36 | ug/m3 |
| 80-62-6 | 100.12 | 5 | ND | 0.80 | 0.12 | ppbv | ND | 3.3 | 0.49 | ug/m3 |
| 115-07-1 | 42 | Propylene | 2.3 | 2.0 | 0.32 | ppbv | 4.0 | 3.4 | 0.55 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.10 | ppbv | ND | 3.4 | 0.43 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.13 | ppbv | ND | 4.4 | 0.71 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.12 | ppbv | ND | 5.5 | 0.82 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.14 | ppbv | ND | 4.4 | 0.76 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.18 | ppbv | ND | 5.9 | 1.3 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 4.3 | 0.80 | 0.092 | ppbv | 21 | 3.9 | 0.45 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 1.4 | 0.80 | 0.12 | ppbv | 6.9 | 3.9 | 0.59 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 0.55 | 0.80 | 0.085 | ppbv J | 2.6 | 3.7 | 0.40 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 9.5 | 0.80 | 0.20 | ppbv | 29 | 2.4 | 0.61 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 2.4 | 0.16 | 0.095 | ppbv | 16 | 1.1 | 0.64 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.80 | 0.17 | ppbv | ND | 2.4 | 0.50 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 17.9 | 0.80 | 0.081 | ppbv | 67.5 | 3.0 | 0.31 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | 0.16 | 0.10 | ppbv | ND | 0.86 | 0.54 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.081 | ppbv | ND | 4.5 | 0.46 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.13 | ppbv | ND | 2.0 | 0.33 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 15.9 | 0.80 | 0.17 | ppbv | 69.1 | 3.5 | 0.74 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 5.5 | 0.80 | 0.10 | ppbv | 24 | 3.5 | 0.43 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 21.5 | 0.80 | 0.10 | ppbv | 93.4 | 3.5 | 0.43 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | ± 2 | Limits | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 104% | 97% | | 65-128% |) | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Page 2 of 3

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3.2

13 of 23

| | Report of Analysis | Page 3 of 3 | 3.2 |
|---|---|-------------|-----|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-7 JC7667-2 Date Sampled: 10/30/15 AIR - Soil Vapor Comp. Summa ID: A873 Date Received: 11/03/15 TO-15 Percent Solids: n/a 2002-2024 Cropsey Avenue, Brooklyn, NY n/a | | ယ |
| VOA TO15 List | | | |
| CAS No. MW | Compound Result RL MDL Units Q Result RL | MDL Uni | its |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



14 of 23 ACCUTEST JC7667

| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC766 AIR - TO-15 | Soil Vapor | Comp. Summ ey Avenue, Bro | | '9 I F | Date Received: 11 | 10/30/15 11/03/15 n/a | |
|---|--|------------|------------------------------|-----|-----------|-------------------|-----------------------------|--|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| Run #1 | 3W50846.D | 1.48 | 11/05/15 | YMH | n/a | n/a | V3W1931 | |
| Run #2 | 3W50860.D | 1.48 | 11/06/15 | YMH | n/a | n/a | V3W1932 | |
| Run #1 Run #2 | Initial Volume 148 ml 50.0 ml | ; | | | | | | |

Report of Analysis

Page 1 of 3

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VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|-------|---|------------------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 267 ^a | 2.4 | 0.38 | ppbv | | 634 ^a | 5.7 | 0.90 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.12 | ppbv | | ND | 1.8 | 0.27 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 1.5 | 0.80 | 0.12 | ppbv | | 4.8 | 2.6 | 0.38 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.13 | ppbv | | ND | 5.4 | 0.87 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.082 | ppbv | | ND | 8.3 | 0.85 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.087 | ppbv | | ND | 3.1 | 0.34 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.081 | ppbv | | ND | 3.5 | 0.35 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.10 | ppbv | | ND | 4.1 | 0.52 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.13 | ppbv | | ND | 3.7 | 0.60 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.086 | ppbv | | ND | 2.1 | 0.23 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | 8.2 | 0.80 | 0.12 | ppbv | | 40 | 3.9 | 0.59 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.12 | ppbv | | ND | 1.7 | 0.25 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.13 | ppbv | | ND | 4.1 | 0.67 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.098 | ppbv | | ND | 5.0 | 0.62 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 0.62 | 0.80 | 0.13 | ppbv | J | 2.1 | 2.8 | 0.45 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.12 | ppbv | | ND | 3.2 | 0.49 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | | ND | 3.2 | 0.44 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.14 | ppbv | | ND | 6.1 | 1.1 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.10 | ppbv | | ND | 3.2 | 0.40 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.20 | ppbv | | ND | 3.7 | 0.92 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.25 | ppbv | | ND | 2.9 | 0.90 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.57 | 0.80 | 0.15 | ppbv | J | 2.8 | 4.0 | 0.74 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.17 | ppbv | | ND | 6.8 | 1.4 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.081 | ppbv | | ND | 3.2 | 0.32 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.099 | ppbv | | ND | 3.2 | 0.39 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.14 | ppbv | | ND | 3.6 | 0.64 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | | ND | 4.8 | 0.66 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.12 | ppbv | | ND | 4.8 | 0.72 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.077 | ppbv | | ND | 4.8 | 0.46 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.082 | ppbv | | ND | 3.6 | 0.37 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



| Client Sample ID: | SV-6 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7667-3 | Date Sampled: | 10/30/15 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A779 | Date Received: | 11/03/15 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|------|---------|--------|---|--------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 7.9 | 2.0 | 0.66 | ppbv | | 15 | 3.8 | 1.2 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 4.7 | 0.80 | 0.19 | ppbv | | 20 | 3.5 | 0.83 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | ND | 0.80 | 0.25 | ppbv | | ND | 2.9 | 0.90 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | 1.9 | 0.80 | 0.088 | ppbv | | 9.3 | 3.9 | 0.43 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.11 | ppbv | | ND | 6.1 | 0.84 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.10 | ppbv | | ND | 5.6 | 0.70 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 4.7 | 0.80 | 0.12 | ppbv | | 19 | 3.3 | 0.49 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.13 | ppbv | | ND | 8.5 | 1.4 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 1.6 | 0.80 | 0.11 | ppbv | | 5.6 | 2.8 | 0.39 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | 6.5 | 0.80 | 0.18 | ppbv | | 27 | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | ND | 0.80 | 0.48 | ppbv | | ND | 2.0 | 1.2 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.54 | ppbv | | ND | 2.8 | 1.9 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 19.3 | 0.80 | 0.19 | ppbv | | 56.9 | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.80 | 0.11 | ppbv | | ND | 3.3 | 0.45 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.10 | ppbv | | ND | 2.9 | 0.36 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.12 | ppbv | | ND | 3.3 | 0.49 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 2.0 | 0.32 | ppbv | | ND | 3.4 | 0.55 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.10 | ppbv | | ND | 3.4 | 0.43 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.13 | ppbv | | ND | 4.4 | 0.71 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.12 | ppbv | | ND | 5.5 | 0.82 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.14 | ppbv | | ND | 4.4 | 0.76 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.18 | ppbv | | ND | 5.9 | 1.3 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 5.8 | 0.80 | 0.092 | ppbv | | 29 | 3.9 | 0.45 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 1.9 | 0.80 | 0.12 | ppbv | | 9.3 | 3.9 | 0.59 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 0.57 | 0.80 | 0.085 | 11 | J | 2.7 | 3.7 | 0.40 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 7.3 | 0.80 | 0.20 | ppbv | | 22 | 2.4 | 0.61 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 111 | 0.16 | 0.095 | ppbv | | 753 | 1.1 | 0.64 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.80 | 0.17 | ppbv | | ND | 2.4 | 0.50 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 18.4 | 0.80 | 0.081 | ppbv | | 69.3 | 3.0 | 0.31 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 0.25 | 0.16 | 0.10 | ppbv | _ | 1.3 | 0.86 | 0.54 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | 0.32 | 0.80 | 0.081 | 11 | J | 1.8 | 4.5 | 0.46 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.13 | ppbv | | ND | 2.0 | 0.33 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 18.6 | 0.80 | 0.17 | ppbv | | 80.8 | 3.5 | 0.74 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 6.2 | 0.80 | 0.10 | ppbv | | 27 | 3.5 | 0.43 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 24.8 | 0.80 | 0.10 | ppbv | | 108 | 3.5 | 0.43 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | l Run# | 2 | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 104% | 101% | | 65-128% | ,) | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Page 2 of 3

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16 of 23

| | Report of Analysis | Page 3 of 3 | ယ ယ |
|---|--|-------------|--------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-6 JC7667-3 Date Sampled: 10/30/15 AIR - Soil Vapor Comp. Summa ID: A779 Date Received: 11/03/15 TO-15 Percent Solids: n/a | | ယ |
| VOA TO15 List | | , | |
| CAS No. MW | Compound Result RL MDL Units Q Result RL | MDL Un | nits |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



17 of 23



Section 4

Sample Results

Report of Analysis





| Client San Lab Samj Matrix: Method: Project: | ple ID: JC155 AIR - TO-15 | 05-1 Soil Vapor | Comp. Summ ey Avenue, Bro | Date Sampled: Date Received: Percent Solids: | | | |
|--|---------------------------------|--------------------|------------------------------|--|-----------|------------|------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | 5W16969.D | 1.28 | 03/13/16 | TCH | n/a | n/a | V5W679 |
| Run #2 | 5W16946.D | 128 | 03/12/16 | ТСН | n/a | n/a | V5W678 |
| | Initial Volum | e | | | | | |
| Run #1 | 20.0 ml | | | | | | |
| Run #2 | 200 ml | | | | | | |

Report of Analysis

Page 1 of 3

4

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|-----|------|---------|--------|----|-----|-------|
| 67-64-1 | 58.08 | Acetone | ND | 5.1 | 0.93 | ppbv | ND | 12 | 2.2 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 5.1 | 0.71 | ppbv | ND | 11 | 1.6 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 27.9 | 5.1 | 0.80 | ppbv | 89.1 | 16 | 2.6 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 5.1 | 0.99 | ppbv | ND | 34 | 6.6 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 5.1 | 0.40 | ppbv | ND | 53 | 4.1 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 5.1 | 0.47 | ppbv | ND | 20 | 1.8 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 5.1 | 0.47 | ppbv | ND | 22 | 2.1 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 5.1 | 0.69 | ppbv | ND | 26 | 3.6 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | 2.5 | 5.1 | 0.80 | ppbv J | 7.8 | 16 | 2.5 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 5.1 | 1.4 | ppbv | ND | 23 | 6.4 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | 5.3 | 5.1 | 0.91 | ppbv | 14 | 13 | 2.4 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 5.1 | 0.42 | ppbv | ND | 25 | 2.1 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 5.1 | 1.3 | ppbv | ND | 11 | 2.7 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 5.1 | 0.68 | ppbv | ND | 16 | 2.1 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 5.1 | 0.43 | ppbv | ND | 26 | 2.2 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 5.1 | 0.80 | ppbv | ND | 32 | 5.0 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 631 | 5.1 | 0.41 | ppbv | 2170 | 18 | 1.4 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 5.1 | 0.39 | ppbv | ND | 21 | 1.6 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 5.1 | 0.54 | ppbv | ND | 20 | 2.1 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 5.1 | 1.1 | ppbv | ND | 39 | 8.5 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 5.1 | 0.45 | ppbv | ND | 21 | 1.8 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 5.1 | 0.56 | ppbv | ND | 24 | 2.6 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 5.1 | 1.2 | ppbv | ND | 18 | 4.3 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | ND | 5.1 | 0.49 | ppbv | ND | 25 | 2.4 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 5.1 | 1.3 | ppbv | ND | 43 | 11 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | 4.5 | 5.1 | 0.72 | ppbv J | 18 | 20 | 2.9 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 5.1 | 0.54 | ppbv | ND | 20 | 2.1 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 5.1 | 0.39 | ppbv | ND | 23 | 1.8 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 5.1 | 0.50 | ppbv | ND | 31 | 3.0 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 5.1 | 0.41 | ppbv | ND | 31 | 2.5 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 5.1 | 0.69 | ppbv | ND | 31 | 4.1 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 5.1 | 0.47 | ppbv | ND | 23 | 2.1 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



| Client Sample ID: | SV-11 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC15505-1 | Date Sampled: | 03/04/16 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A1177,A1216 | Date Received: | 03/05/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|-------------------|------------|---------|---------|-------------------|-----|-----|-------|
| 64-17-5 | 46.07 | Ethanol | 22.2 | 13 | 1.9 | ppbv | 41.8 | 24 | 3.6 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | ND | 5.1 | 1.1 | ppbv | ND | 22 | 4.8 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | ND | 5.1 | 1.9 | ppbv | ND | 18 | 6.8 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 5.1 | 0.43 | ppbv | ND | 25 | 2.1 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 5.1 | 0.55 | ppbv | ND | 39 | 4.2 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 5.1 | 0.80 | ppbv | ND | 36 | 5.6 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 360 | 5.1 | 0.52 | ppbv | 1480 | 21 | 2.1 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 5.1 | 0.52 | ppbv | ND | 54 | 5.5 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 1640 ^a | 51 | 5.8 | ppbv | 5780 ^a | 180 | 20 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 5.1 | 1.1 | ppbv | ND | 21 | 4.5 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | ND | 5.1 | 4.0 | ppbv | ND | 13 | 9.8 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 5.1 | 0.64 | ppbv | ND | 18 | 2.2 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 344 | 5.1 | 1.2 | ppbv | 1010 | 15 | 3.5 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 5.1 | 1.4 | ppbv | ND | 21 | 5.7 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 5.1 | 0.50 | ppbv | ND | 18 | 1.8 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 5.1 | 1.0 | ppbv | ND | 21 | 4.1 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 13 | 0.83 | ppbv | ND | 22 | 1.4 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 5.1 | 0.39 | ppbv | ND | 22 | 1.7 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 5.1 | 0.60 | ppbv | ND | 28 | 3.3 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 5.1 | 0.41 | ppbv | ND | 35 | 2.8 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 5.1 | 0.99 | ppbv | ND | 28 | 5.4 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 5.1 | 1.4 | ppbv | ND | 38 | 10 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | ND | 5.1 | 0.39 | ppbv | ND | 25 | 1.9 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | 5.1 | 1.1 | ppbv | ND | 25 | 5.4 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | ND | 5.1 | 0.58 | ppbv | ND | 24 | 2.7 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 18.6 | 5.1 | 1.4 | ppbv | 56.4 | 15 | 4.2 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 1.1 | 1.0 | 0.59 | ppbv | 7.5 | 6.8 | 4.0 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | 830 | 5.1 | 1.2 | ppbv | 2450 | 15 | 3.5 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | ND | 5.1 | 0.32 | ppbv | ND | 19 | 1.2 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 1.9 | 1.0 | 0.48 | ppbv | 10 | 5.4 | 2.6 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 5.1 | 0.56 | ppbv | ND | 29 | 3.1 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | 8.9 | 5.1 | 0.53 | ppbv | 23 | 13 | 1.4 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 5.1 | 1.4 | ppbv | ND | 18 | 4.9 | ug/m3 |
| | 106.2 | m,p-Xylene | ND | 5.1 | 1.7 | ppbv | ND | 22 | 7.4 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | ND | 5.1 | 1.3 | ppbv | ND | 22 | 5.6 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | ND | 5.1 | 1.3 | ppbv | ND | 22 | 5.6 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | # 2 | Limits | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 100% | 95% | | 65-128% | ,) | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



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12 of 507

ACCUTEST JC15505

| | Report of Analysis | | Page 3 of 3 |
|---|---|--|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-11 JC15505-1 AIR - Soil Vapor Comp. Summa ID: A1177,A1216 TO-15 2002-2024 Cropsey Avenue, Brooklyn, NY | Date Sampled: 03/04/16 Date Received: 03/05/16 Percent Solids: n/a | |
| VOA TO15 List | | | |
| CAS No. MW | Compound Result RL MDI | Units Q Result RI | L MDL Units |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



13 of 507

| Lab Samj Matrix: Method: Project: | AIR - TO-1 | 505-2 Ambient A 5 | Air Comp. Sum sey Avenue, Bro | nma ID: A ooklyn, NY | | Date Sampled: Date Received: Percent Solids: | |
|--|---------------|-------------------------|----------------------------------|-------------------------|-----------|--|--------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batc | h Analytical Batch |
| Run #1 | 5W16947.D | 1 | 03/12/16 | TCH | n/a | n/a | V5W678 |
| Run #2 | | | | | | | |
| | Initial Volum | e | | | | | |
| Run #1 | 400 ml | | | | | | |
| D 110 | | | | | | | |

Report of Analysis

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|------|-------|---------|--------|------|-------|-------|
| 67-64-1 | 58.08 | Acetone | 4.1 | 0.20 | 0.036 | ppbv | 9.7 | 0.48 | 0.086 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | 0.25 | 0.20 | 0.028 | | 0.55 | 0.44 | 0.062 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 2.5 | 0.20 | 0.031 | ppbv | 8.0 | 0.64 | 0.099 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.20 | 0.039 | ppbv | ND | 1.3 | 0.26 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.20 | 0.016 | | ND | 2.1 | 0.17 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.20 | 0.018 | ppbv | ND | 0.78 | 0.070 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.20 | 0.018 | ppbv | ND | 0.87 | 0.079 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.20 | 0.027 | ppbv | ND | 1.0 | 0.14 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.20 | 0.031 | ppbv | ND | 0.62 | 0.097 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.20 | 0.056 | ppbv | ND | 0.92 | 0.26 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.20 | 0.036 | | ND | 0.53 | 0.095 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.20 | 0.017 | | ND | 0.98 | 0.083 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | 0.63 | 0.20 | 0.052 | | 1.3 | 0.41 | 0.11 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.20 | 0.027 | | ND | 0.63 | 0.085 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.20 | 0.017 | | ND | 1.0 | 0.088 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.20 | 0.031 | | ND | 1.3 | 0.20 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 0.29 | 0.20 | 0.016 | | 1.0 | 0.69 | 0.055 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.20 | 0.015 | | ND | 0.81 | 0.061 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.20 | 0.021 | | ND | 0.79 | 0.083 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.20 | 0.042 | | ND | 1.5 | 0.32 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.20 | 0.018 | | ND | 0.81 | 0.073 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.20 | 0.022 | | ND | 0.92 | 0.10 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.20 | 0.045 | | ND | 0.72 | 0.16 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.56 | 0.20 | 0.019 | 11 | 2.8 | 0.99 | 0.094 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.20 | 0.053 | 11 | ND | 1.7 | 0.45 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.20 | 0.028 | | ND | 0.79 | 0.11 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.20 | 0.021 | | ND | 0.79 | 0.083 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.20 | 0.015 | | ND | 0.91 | 0.068 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.20 | 0.020 | | ND | 1.2 | 0.12 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.20 | 0.016 | | ND | 1.2 | 0.096 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.20 | 0.027 | 11 | ND | 1.2 | 0.16 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.20 | 0.018 | ppbv | ND | 0.91 | 0.082 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2



14 of 507

| Report | of | Ana | lysis | |
|--------|----|-----|-------|--|
|--------|----|-----|-------|--|

| Client Sample ID: | IAQ-11 | | |
|-------------------|---|------------------------|----------|
| Lab Sample ID: | JC15505-2 | Date Sampled: | 03/04/16 |
| Matrix: | AIR - Ambient Air Comp. Summa ID: A1112 | Date Received: | 03/05/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

VOA TO15 List

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| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|-------|---------|------------|---|--------|------|-------|-------|
| 64-17-5 | 46.07 | Ethanol | 15.9 | 0.50 | 0.075 | ppbv | | 30.0 | 0.94 | 0.14 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 0.32 | 0.20 | 0.042 | ppbv | | 1.4 | 0.87 | 0.18 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 1.2 | 0.20 | 0.075 | | | 4.3 | 0.72 | 0.27 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 0.20 | 0.017 | ppbv | | ND | 0.98 | 0.084 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.20 | 0.021 | ppbv | | ND | 1.5 | 0.16 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.20 | 0.031 | ppbv | | ND | 1.4 | 0.22 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 0.35 | 0.20 | 0.020 | | | 1.4 | 0.82 | 0.082 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.20 | 0.020 | | | ND | 2.1 | 0.21 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 0.81 | 0.20 | 0.023 | ppbv | | 2.9 | 0.70 | 0.081 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 0.20 | 0.045 | ppbv | | ND | 0.82 | 0.18 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 7.2 | 0.20 | 0.16 | ppbv | | 18 | 0.49 | 0.39 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | 0.22 | 0.20 | 0.025 | ppbv | | 0.76 | 0.69 | 0.087 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 0.26 | 0.20 | 0.048 | ppbv | | 0.77 | 0.59 | 0.14 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.20 | 0.055 | | | ND | 0.82 | 0.23 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.20 | 0.020 | | | ND | 0.72 | 0.072 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.20 | 0.040 | | | ND | 0.82 | 0.16 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 0.50 | 0.032 | ppbv | | ND | 0.86 | 0.055 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.20 | 0.015 | | | ND | 0.85 | 0.064 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.20 | 0.024 | | | ND | 1.1 | 0.13 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.20 | 0.016 | . . | | ND | 1.4 | 0.11 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.20 | 0.039 | | | ND | 1.1 | 0.21 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.20 | 0.056 | | | ND | 1.5 | 0.42 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 0.33 | 0.20 | 0.015 | | | 1.6 | 0.98 | 0.074 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 0.11 | 0.20 | 0.045 | ppbv | J | 0.54 | 0.98 | 0.22 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 1.1 | 0.20 | 0.023 | | | 5.1 | 0.93 | 0.11 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | ND | 0.20 | 0.053 | | | ND | 0.61 | 0.16 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 0.072 | 0.040 | 0.023 | ppbv | | 0.49 | 0.27 | 0.16 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.20 | | | | ND | 0.59 | 0.13 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 3.1 | 0.20 | 0.012 | | | 12 | 0.75 | 0.045 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | 0.040 | | | | ND | 0.21 | 0.10 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | 0.25 | 0.20 | 0.022 | | | 1.4 | 1.1 | 0.12 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.20 | 0.021 | | | ND | 0.51 | 0.054 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.20 | 0.054 | | | ND | 0.70 | 0.19 | ug/m3 |
| | 106.2 | m,p-Xylene | 1.4 | 0.20 | 0.068 | | | 6.1 | 0.87 | 0.30 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 0.49 | 0.20 | 0.051 | | | 2.1 | 0.87 | 0.22 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 1.9 | 0.20 | 0.051 | ppbv | | 8.3 | 0.87 | 0.22 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run | #2 I | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 103% | | e | 55-128% |) | | | | | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



4.2

4

Page 2 of 2



| | | | Repo | ort of A | nalysis | | Page 1 of 2 |
|--|-------------------------------|-----------------------------|--------------------------|-------------------------|-------------------------|--|------------------------------|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC1 AIR TO- | 5505-3 - Ambient A 15 | Air Comp. Sum | nma ID: A poklyn, NY | | Date Sampled: Date Received: Percent Solids: | |
| Run #1 Run #2 | File ID 5W16948.D | DF 1 | Analyzed 03/13/16 | By TCH | Prep Date n/a | Prep Batc n/a | h Analytical Batch V5W678 |
| Run #1 | Initial Volu 400 ml | me | | | | | |

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|------|-------|---------|--------|------|-------|-------|
| 67-64-1 | 58.08 | Acetone | 8.7 | 0.20 | 0.036 | ppbv | 21 | 0.48 | 0.086 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | 0.34 | 0.20 | 0.028 | ppbv | 0.75 | 0.44 | 0.062 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 0.74 | 0.20 | 0.031 | ppbv | 2.4 | 0.64 | 0.099 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.20 | 0.039 | | ND | 1.3 | 0.26 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.20 | 0.016 | | ND | 2.1 | 0.17 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.20 | 0.018 | ppbv | ND | 0.78 | 0.070 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.20 | 0.018 | ppbv | ND | 0.87 | 0.079 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.20 | 0.027 | ppbv | ND | 1.0 | 0.14 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.20 | 0.031 | ppbv | ND | 0.62 | 0.097 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.20 | 0.056 | ppbv | ND | 0.92 | 0.26 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.20 | 0.036 | ppbv | ND | 0.53 | 0.095 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.20 | 0.017 | ppbv | ND | 0.98 | 0.083 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | 0.59 | 0.20 | 0.052 | ppbv | 1.2 | 0.41 | 0.11 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.20 | 0.027 | ppbv | ND | 0.63 | 0.085 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.20 | 0.017 | ppbv | ND | 1.0 | 0.088 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.20 | 0.031 | ppbv | ND | 1.3 | 0.20 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 0.13 | 0.20 | 0.016 | ppbv J | 0.45 | 0.69 | 0.055 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.20 | 0.015 | ppbv | ND | 0.81 | 0.061 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.20 | 0.021 | ppbv | ND | 0.79 | 0.083 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.20 | 0.042 | ppbv | ND | 1.5 | 0.32 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.20 | 0.018 | ppbv | ND | 0.81 | 0.073 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.20 | 0.022 | | ND | 0.92 | 0.10 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.20 | 0.045 | ppbv | ND | 0.72 | 0.16 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.49 | 0.20 | 0.019 | ppbv | 2.4 | 0.99 | 0.094 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.20 | 0.053 | ppbv | ND | 1.7 | 0.45 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.20 | 0.028 | ppbv | ND | 0.79 | 0.11 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.20 | 0.021 | ppbv | ND | 0.79 | 0.083 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.20 | 0.015 | | ND | 0.91 | 0.068 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.20 | 0.020 | | ND | 1.2 | 0.12 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.20 | 0.016 | | ND | 1.2 | 0.096 | - |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.20 | 0.027 | | ND | 1.2 | 0.16 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.20 | 0.018 | | ND | 0.91 | 0.082 | - |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



JC15505

4.3

| Report of Analysis |
|---------------------------|
| |

| Client Sample ID: | | | |
|-------------------|---|-----------------|----------|
| Lab Sample ID: | JC15505-3 | Date Sampled: | 03/04/16 |
| Matrix: | AIR - Ambient Air Comp. Summa ID: A1183 | Date Received: | 03/05/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|-------|---------|-------|---|--------|------|-------|-------|
| 64-17-5 | 46.07 | Ethanol | 8.9 | 0.50 | 0.075 | ppbv | | 17 | 0.94 | 0.14 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 0.18 | 0.20 | 0.042 | | J | 0.78 | 0.87 | 0.18 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 1.6 | 0.20 | 0.075 | ppbv | | 5.8 | 0.72 | 0.27 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 0.20 | 0.017 | ppbv | | ND | 0.98 | 0.084 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.20 | 0.021 | | | ND | 1.5 | 0.16 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.20 | 0.031 | ppbv | | ND | 1.4 | 0.22 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 0.17 | 0.20 | 0.020 | ppbv | J | 0.70 | 0.82 | 0.082 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.20 | 0.020 | ppbv | | ND | 2.1 | 0.21 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 0.60 | 0.20 | 0.023 | ppbv | | 2.1 | 0.70 | 0.081 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 0.20 | 0.045 | ppbv | | ND | 0.82 | 0.18 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 6.7 | 0.20 | 0.16 | ppbv | | 16 | 0.49 | 0.39 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | 0.46 | 0.20 | 0.025 | ppbv | | 1.6 | 0.69 | 0.087 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 0.28 | 0.20 | 0.048 | ppbv | | 0.83 | 0.59 | 0.14 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.20 | 0.055 | ppbv | | ND | 0.82 | 0.23 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.20 | 0.020 | ppbv | | ND | 0.72 | 0.072 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.20 | 0.040 | ppbv | | ND | 0.82 | 0.16 | ug/m3 |
| 115-07-1 | 42 | Propylene | 3.1 | 0.50 | 0.032 | ppbv | | 5.3 | 0.86 | 0.055 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.20 | 0.015 | ppbv | | ND | 0.85 | 0.064 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.20 | 0.024 | ppbv | | ND | 1.1 | 0.13 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.20 | 0.016 | ppbv | | ND | 1.4 | 0.11 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.20 | 0.039 | | | ND | 1.1 | 0.21 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.20 | 0.056 | ppbv | | ND | 1.5 | 0.42 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 0.12 | 0.20 | 0.015 | ppbv | J | 0.59 | 0.98 | 0.074 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | 0.20 | 0.045 | ppbv | | ND | 0.98 | 0.22 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 0.24 | 0.20 | 0.023 | ppbv | | 1.1 | 0.93 | 0.11 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | ND | 0.20 | 0.053 | ppbv | | ND | 0.61 | 0.16 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 0.065 | 0.040 | 0.023 | ppbv | | 0.44 | 0.27 | 0.16 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | 0.23 | 0.20 | 0.045 | ppbv | | 0.68 | 0.59 | 0.13 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 1.2 | 0.20 | 0.012 | ppbv | | 4.5 | 0.75 | 0.045 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | 0.040 | | | | ND | 0.21 | 0.10 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | 0.26 | 0.20 | 0.022 | ppbv | | 1.5 | 1.1 | 0.12 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.20 | 0.021 | ppbv | | ND | 0.51 | 0.054 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.20 | 0.054 | ppbv | | ND | 0.70 | 0.19 | ug/m3 |
| | 106.2 | m,p-Xylene | 0.56 | 0.20 | 0.068 | ppbv | | 2.4 | 0.87 | 0.30 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 0.19 | 0.20 | 0.051 | ppbv | J | 0.83 | 0.87 | 0.22 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 0.76 | 0.20 | 0.051 | ppbv | | 3.3 | 0.87 | 0.22 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run | #2 I | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 104% | | e | 55-128% |) | | | | | |

ND = Not detectedMDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Page 2 of 2

4.3 4



17 of 507

| Lab Sample ID: Aatrix: Aethod: Project: | TO-15 | oor Comp. Sumn opsey Avenue, Bro | | 13 D P | Date Sampled:03Date Received:03Percent Solids:n/ | |
|--|--------|-------------------------------------|-----|-----------|--|------------------|
| File II | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 5W169 | 49.D 1 | 03/13/16 | TCH | n/a | n/a | V5W678 |
| Run #2 | | | | | | |
| Initial | Volume | | | | | |
| Run #1 100 m | | | | | | |

Report of Analysis

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|------|-------|-------|---|--------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 17.3 | 0.80 | 0.14 | ppbv | | 41.1 | 1.9 | 0.33 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | 0.78 | 0.80 | 0.11 | ppbv | J | 1.7 | 1.8 | 0.24 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 2.0 | 0.80 | 0.13 | ppbv | | 6.4 | 2.6 | 0.42 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.15 | ppbv | | ND | 5.4 | 1.0 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.063 | ppbv | | ND | 8.3 | 0.65 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.074 | ppbv | | ND | 3.1 | 0.29 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.074 | ppbv | | ND | 3.5 | 0.32 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.11 | ppbv | | ND | 4.1 | 0.57 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.80 | 0.13 | ppbv | | ND | 2.5 | 0.40 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.22 | ppbv | | ND | 3.7 | 1.0 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.14 | ppbv | | ND | 2.1 | 0.37 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.80 | 0.066 | ppbv | | ND | 3.9 | 0.32 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | 0.67 | 0.80 | 0.21 | ppbv | J | 1.4 | 1.7 | 0.43 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.068 | ppbv | | ND | 4.1 | 0.35 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.12 | ppbv | | ND | 5.0 | 0.75 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | ND | 0.80 | 0.065 | ppbv | | ND | 2.8 | 0.22 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.061 | ppbv | | ND | 3.2 | 0.25 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.084 | ppbv | | ND | 3.2 | 0.33 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.17 | ppbv | | ND | 6.1 | 1.3 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.070 | ppbv | | ND | 3.2 | 0.28 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.088 | ppbv | | ND | 3.7 | 0.41 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.18 | ppbv | | ND | 2.9 | 0.65 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.48 | 0.80 | 0.076 | ppbv | J | 2.4 | 4.0 | 0.38 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.21 | ppbv | | ND | 6.8 | 1.8 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | | ND | 3.2 | 0.44 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.084 | ppbv | | ND | 3.2 | 0.33 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.061 | ppbv | | ND | 3.6 | 0.28 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.078 | ppbv | | ND | 4.8 | 0.47 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.064 | ppbv | | ND | 4.8 | 0.38 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | | ND | 4.8 | 0.66 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.073 | ppbv | | ND | 3.6 | 0.33 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

18 of 507

ACCUTEST JC15505

E = Indicates value exceeds calibration range

| Report of | Analysis | |
|-----------|----------|--|
|-----------|----------|--|

| Client Sample ID: | SV-9 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC15505-4 | Date Sampled: | 03/04/16 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A313 | Date Received: | 03/05/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | * • • | | |

VOA TO15 List

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| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|------------|---------|---------|--------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 12.7 | 2.0 | 0.30 | ppbv | 23.9 | 3.8 | 0.57 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | ND | 0.80 | 0.17 | ppbv | ND | 3.5 | 0.74 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 1.5 | 0.80 | 0.30 | ppbv | 5.4 | 2.9 | 1.1 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 0.80 | 0.067 | ppbv | ND | 3.9 | 0.33 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.086 | ppbv | ND | 6.1 | 0.66 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.13 | ppbv | ND | 5.6 | 0.91 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | ND | 0.80 | 0.081 | ppbv | ND | 3.3 | 0.33 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.081 | ppbv | ND | 8.5 | 0.86 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 0.84 | 0.80 | 0.090 | ppbv | 3.0 | 2.8 | 0.32 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 0.80 | 0.18 | ppbv | ND | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 13.9 | 0.80 | 0.62 | ppbv | 34.2 | 2.0 | 1.5 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.10 | ppbv | ND | 2.8 | 0.35 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | ND | 0.80 | 0.19 | ppbv | ND | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.80 | 0.22 | ppbv | ND | 3.3 | 0.90 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.079 | ppbv | ND | 2.9 | 0.28 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.16 | ppbv | ND | 3.3 | 0.66 | ug/m3 |
| 115-07-1 | 42 | Propylene | 4.6 | 2.0 | 0.13 | ppbv | 7.9 | 3.4 | 0.22 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.062 | ppbv | ND | 3.4 | 0.26 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.094 | ppbv | ND | 4.4 | 0.51 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.064 | ppbv | ND | 5.5 | 0.44 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.16 | ppbv | ND | 4.4 | 0.87 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.22 | ppbv | ND | 5.9 | 1.6 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | ND | 0.80 | 0.061 | ppbv | ND | 3.9 | 0.30 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | 0.80 | 0.18 | ppbv | ND | 3.9 | 0.88 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 0.47 | 0.80 | 0.091 | ppbv J | 2.2 | 3.7 | 0.43 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | ND | 0.80 | 0.21 | ppbv | ND | 2.4 | 0.64 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 0.16 | 0.16 | 0.092 | ppbv | 1.1 | 1.1 | 0.62 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.80 | 0.18 | ppbv | ND | 2.4 | 0.53 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 2.1 | 0.80 | 0.050 | ppbv | 7.9 | 3.0 | 0.19 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | 0.16 | 0.074 | ppbv | ND | 0.86 | 0.40 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.088 | ppbv | ND | 4.5 | 0.49 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.082 | ppbv | ND | 2.0 | 0.21 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 0.80 | 0.80 | 0.27 | ppbv | 3.5 | 3.5 | 1.2 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | ND | 0.80 | 0.20 | ppbv | ND | 3.5 | 0.87 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 0.80 | 0.80 | 0.20 | ppbv | 3.5 | 3.5 | 0.87 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | # 2 | Limits | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 101% | | (| 65-128% | ,) | | | | |

ND = Not detectedMDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 2



19 of 507

| Client San Lab Samj Matrix: Method: Project: | AIR - TO-1 | 505-5 · Ambient A 5 | ir Comp. Sum ey Avenue, Bro | Date Sampled: Date Received: Percent Solids: | | | |
|--|---|---------------------------|--------------------------------|--|-----------|-----------|--------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batc | h Analytical Batch |
| Run #1 | 5W16950.D | 1 | 03/13/16 | TCH | n/a | n/a | V5W678 |
| Run #2 | 5W16970.D | 1.55 | 03/13/16 | TCH | n/a | n/a | V5W679 |
| Run #1 Run #2 | Initial Volum 400 ml 20.0 ml | e | | | | | |

Report of Analysis

Page 1 of 3

4.5

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|------------|------------------|------|-------|-------|
| 67-64-1 | 58.08 | Acetone | 153 ^a | 6.2 | 1.1 | ppbv | 363 ^a | 15 | 2.6 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.20 | 0.028 | ppbv | ND | 0.44 | 0.062 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 0.20 | 0.20 | 0.031 | ppbv | 0.64 | 0.64 | 0.099 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.20 | 0.039 | ppbv | ND | 1.3 | 0.26 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.20 | 0.016 | ppbv | ND | 2.1 | 0.17 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.20 | 0.018 | ppbv | ND | 0.78 | 0.070 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.20 | 0.018 | ppbv | ND | 0.87 | 0.079 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.20 | 0.027 | ppbv | ND | 1.0 | 0.14 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | 0.70 | 0.20 | 0.031 | ppbv | 2.2 | 0.62 | 0.097 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.20 | 0.056 | ppbv | ND | 0.92 | 0.26 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.20 | 0.036 | | ND | 0.53 | 0.095 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | 0.23 | 0.20 | 0.017 | ppbv | 1.1 | 0.98 | 0.083 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.20 | 0.052 | ppbv | ND | 0.41 | 0.11 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.20 | 0.027 | ppbv | ND | 0.63 | 0.085 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.20 | 0.017 | ppbv | ND | 1.0 | 0.088 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.20 | 0.031 | ppbv | ND | 1.3 | 0.20 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 0.28 | 0.20 | 0.016 | ppbv | 0.96 | 0.69 | 0.055 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.20 | 0.015 | | ND | 0.81 | 0.061 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.20 | 0.021 | T T | ND | 0.79 | 0.083 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.20 | 0.042 | ppbv | ND | 1.5 | 0.32 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.20 | 0.018 | ppbv | ND | 0.81 | 0.073 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.20 | 0.022 | ppbv | ND | 0.92 | 0.10 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.20 | 0.045 | * * | ND | 0.72 | 0.16 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.54 | 0.20 | 0.019 | ppbv | 2.7 | 0.99 | 0.094 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.20 | 0.053 | T T | ND | 1.7 | 0.45 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.20 | 0.028 | ppbv | ND | 0.79 | 0.11 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.20 | 0.021 | ppbv | ND | 0.79 | 0.083 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.20 | 0.015 | ppbv | ND | 0.91 | 0.068 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.20 | 0.020 | 11 | ND | 1.2 | 0.12 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.20 | 0.016 | | ND | 1.2 | 0.096 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.20 | 0.027 | | ND | 1.2 | 0.16 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.20 | 0.018 | ppbv | ND | 0.91 | 0.082 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



| Client Sample ID: | IAQ-9 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC15505-5 | Date Sampled: | 03/04/16 |
| Matrix: | AIR - Ambient Air Comp. Summa ID: A278 | Date Received: | 03/05/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|------------------|-------|--------|------------|---|-------------------|------|-------|-------|
| 64-17-5 | 46.07 | Ethanol | 15.5 | 0.50 | 0.075 | ppbv | | 29.2 | 0.94 | 0.14 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 0.11 | 0.20 | 0.042 | ppbv | J | 0.48 | 0.87 | 0.18 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 2.0 | 0.20 | 0.075 | ppbv | | 7.2 | 0.72 | 0.27 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 0.20 | 0.017 | ppbv | | ND | 0.98 | 0.084 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.20 | 0.021 | | | ND | 1.5 | 0.16 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.20 | 0.031 | ppbv | | ND | 1.4 | 0.22 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 0.12 | 0.20 | 0.020 | ppbv | J | 0.49 | 0.82 | 0.082 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.20 | 0.020 | ppbv | | ND | 2.1 | 0.21 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 0.26 | 0.20 | 0.023 | ppbv | | 0.92 | 0.70 | 0.081 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | 0.60 | 0.20 | 0.045 | ppbv | | 2.5 | 0.82 | 0.18 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 8.2 | 0.20 | 0.16 | ppbv | | 20 | 0.49 | 0.39 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | 0.18 | 0.20 | 0.025 | ppbv | J | 0.63 | 0.69 | 0.087 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 133 a | 6.2 | 1.5 | ppbv | | 392 a | 18 | 4.4 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | 1.2 | 0.20 | 0.055 | ppbv | | 4.9 | 0.82 | 0.23 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.20 | 0.020 | ppbv | | ND | 0.72 | 0.072 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.20 | 0.040 | ppbv | | ND | 0.82 | 0.16 | ug/m3 |
| 115-07-1 | 42 | Propylene | 1.5 | 0.50 | 0.032 | ppbv | | 2.6 | 0.86 | 0.055 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.20 | 0.015 | ppbv | | ND | 0.85 | 0.064 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.20 | 0.024 | ppbv | | ND | 1.1 | 0.13 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.20 | 0.016 | | | ND | 1.4 | 0.11 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.20 | 0.039 | ppbv | | ND | 1.1 | 0.21 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.20 | 0.056 | ppbv | | ND | 1.5 | 0.42 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 0.32 | 0.20 | 0.015 | ppbv | | 1.6 | 0.98 | 0.074 | U |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 0.097 | 0.20 | 0.045 | ppbv | J | 0.48 | 0.98 | 0.22 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | ND | 0.20 | 0.023 | ppbv | | ND | 0.93 | 0.11 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 6.5 | 0.20 | 0.053 | ppbv | | 20 | 0.61 | 0.16 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 216 ^a | 1.2 | 0.71 | ppbv | | 1460 ^a | 8.1 | 4.8 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | 397 a | 6.2 | 1.4 | ppbv | | 1170 a | 18 | 4.1 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 0.90 | 0.20 | 0.012 | | | 3.4 | 0.75 | 0.045 | U |
| 79-01-6 | 131.4 | Trichloroethylene | 0.19 | 0.040 | 0.019 | ppbv | | 1.0 | 0.21 | 0.10 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | 0.26 | 0.20 | 0.022 | | | 1.5 | 1.1 | 0.12 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.20 | 0.021 | | | ND | 0.51 | 0.054 | U |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.20 | 0.054 | | | ND | 0.70 | 0.19 | ug/m3 |
| | 106.2 | m,p-Xylene | 0.45 | 0.20 | 0.068 | . . | | 2.0 | 0.87 | 0.30 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 0.29 | 0.20 | 0.051 | | | 1.3 | 0.87 | 0.22 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 0.73 | 0.20 | 0.051 | ppbv | | 3.2 | 0.87 | 0.22 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | 2 I | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 105% | 101% | e | 5-128% |) | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



JC15505

Page 2 of 3

4.5 4

| | Report of Analys | sis | Page 3 of 3 |
|---|---|---|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | IAQ-9 JC15505-5 AIR - Ambient Air Comp. Summa ID: A278 TO-15 2002-2024 Cropsey Avenue, Brooklyn, NY | Date Sampled:03/04/16Date Received:03/05/16Percent Solids:n/a | 4.5 4 |
| VOA TO15 List | | | |
| CAS No. MW | Compound Result RL | MDL Units Q Result RL | MDL Units |

(a) Result is from Run# 2

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



22 of 507 ACCUTEST JC15505



Section 4

Sample Results

Report of Analysis





| Client Sar Lab Samj Matrix: Method: Project: | ple ID: JC15 AIR - TO-1 | Ambient A 5 | Air Comp. Sum sey Avenue, Bro | nma ID: A ooklyn, NY | | Date Sampled: Date Received: Percent Solids: | |
|--|--------------------------------|----------------|----------------------------------|-------------------------|-------------------------|--|------------------------------|
| Run #1 Run #2 | File ID 5W17095.D | DF 1 | Analyzed 03/18/16 | By TCH | Prep Date n/a | Prep Batcl n/a | h Analytical Batch V5W683 |
| Run #1 | Initial Volum 400 ml | e | | | | | |

Report of Analysis

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|------|-------|-----------|--------|------|-------|-------|
| 67-64-1 | 58.08 | Acetone | 5.5 | 0.20 | 0.036 | ppbv | 13 | 0.48 | 0.086 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.20 | 0.028 | | ND | 0.44 | 0.062 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 0.33 | 0.20 | 0.031 | ppbv | 1.1 | 0.64 | 0.099 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.20 | 0.039 | ppbv | ND | 1.3 | 0.26 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.20 | 0.016 | | ND | 2.1 | 0.17 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.20 | 0.018 | ppbv | ND | 0.78 | 0.070 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.20 | 0.018 | | ND | 0.87 | 0.079 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.20 | 0.027 | ppbv | ND | 1.0 | 0.14 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.20 | 0.031 | ppbv | ND | 0.62 | 0.097 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.20 | 0.056 | * * | ND | 0.92 | 0.26 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.20 | 0.036 | | ND | 0.53 | 0.095 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.20 | 0.017 | | ND | 0.98 | 0.083 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | 0.87 | 0.20 | 0.052 | | 1.8 | 0.41 | 0.11 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.20 | 0.027 | | ND | 0.63 | 0.085 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.20 | 0.017 | | ND | 1.0 | 0.088 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | 0.096 | 0.20 | | ppbv J | 0.60 | 1.3 | 0.20 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | ND | 0.20 | 0.016 | | ND | 0.69 | 0.055 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.20 | 0.015 | | ND | 0.81 | 0.061 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.20 | 0.021 | | ND | 0.79 | 0.083 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.20 | 0.042 | | ND | 1.5 | 0.32 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.20 | 0.018 | | ND | 0.81 | 0.073 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.20 | 0.022 | | ND | 0.92 | 0.10 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.20 | 0.045 | | ND | 0.72 | 0.16 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.69 | 0.20 | 0.019 | ~ ~ | 3.4 | 0.99 | 0.094 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.20 | 0.053 | 11 | ND | 1.7 | 0.45 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.20 | 0.028 | | ND | 0.79 | 0.11 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.20 | 0.021 | ppbv | ND | 0.79 | 0.083 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.20 | 0.015 | | ND | 0.91 | 0.068 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.20 | 0.020 | | ND | 1.2 | 0.12 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.20 | 0.016 | | ND | 1.2 | 0.096 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.20 | 0.027 | | ND | 1.2 | 0.16 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.20 | 0.018 | ppbv | ND | 0.91 | 0.082 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

4

Page 1 of 2



| Client Sample ID: Lab Sample ID: | AMBIENT JC15931-1 | Date Sampled: | 03/09/16 |
|-------------------------------------|---|-----------------|----------|
| Matrix: | AIR - Ambient Air Comp. Summa ID: A1071 | Date Received: | |
| Method: Project: | TO-15 2002-2024 Cropsey Avenue, Brooklyn, NY | Percent Solids: | n/a |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|-------|---------|-------|---|--------|------|-------|-------|
| 64-17-5 | 46.07 | Ethanol | 11.5 | 0.50 | 0.075 | ppbv | | 21.7 | 0.94 | 0.14 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | ND | 0.20 | 0.042 | | | ND | 0.87 | 0.18 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 2.6 | 0.20 | 0.075 | ppbv | | 9.4 | 0.72 | 0.27 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 0.20 | 0.017 | ppbv | | ND | 0.98 | 0.084 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | 0.099 | 0.20 | 0.021 | ppbv | J | 0.76 | 1.5 | 0.16 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.20 | 0.031 | ppbv | | ND | 1.4 | 0.22 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 0.10 | 0.20 | 0.020 | ppbv | J | 0.41 | 0.82 | 0.082 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.20 | 0.020 | ppbv | | ND | 2.1 | 0.21 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 0.31 | 0.20 | 0.023 | ppbv | | 1.1 | 0.70 | 0.081 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 0.20 | 0.045 | ppbv | | ND | 0.82 | 0.18 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 2.0 | 0.20 | 0.16 | ppbv | | 4.9 | 0.49 | 0.39 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | 0.49 | 0.20 | 0.025 | ppbv | | 1.7 | 0.69 | 0.087 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 0.43 | 0.20 | 0.048 | ppbv | | 1.3 | 0.59 | 0.14 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 0.20 | 0.055 | ppbv | | ND | 0.82 | 0.23 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.20 | 0.020 | ppbv | | ND | 0.72 | 0.072 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.20 | 0.040 | ppbv | | ND | 0.82 | 0.16 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 0.50 | 0.032 | ppbv | | ND | 0.86 | 0.055 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.20 | 0.015 | ppbv | | ND | 0.85 | 0.064 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.20 | 0.024 | ppbv | | ND | 1.1 | 0.13 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.20 | 0.016 | ppbv | | ND | 1.4 | 0.11 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.20 | 0.039 | ppbv | | ND | 1.1 | 0.21 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.20 | 0.056 | ppbv | | ND | 1.5 | 0.42 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | ND | 0.20 | 0.015 | ppbv | | ND | 0.98 | 0.074 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | 0.20 | 0.045 | ppbv | | ND | 0.98 | 0.22 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 0.18 | 0.20 | 0.023 | ppbv | J | 0.84 | 0.93 | 0.11 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | ND | 0.20 | 0.053 | ppbv | | ND | 0.61 | 0.16 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 0.14 | 0.040 | 0.023 | ppbv | | 0.95 | 0.27 | 0.16 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.20 | 0.045 | ppbv | | ND | 0.59 | 0.13 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 0.75 | 0.20 | 0.012 | ppbv | | 2.8 | 0.75 | 0.045 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | 0.040 | 0.019 | ppbv | | ND | 0.21 | 0.10 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | 0.37 | 0.20 | 0.022 | ppbv | | 2.1 | 1.1 | 0.12 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.20 | 0.021 | ppbv | | ND | 0.51 | 0.054 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.20 | 0.054 | ppbv | | ND | 0.70 | 0.19 | ug/m3 |
| | 106.2 | m,p-Xylene | 0.22 | 0.20 | 0.068 | ppbv | | 0.96 | 0.87 | 0.30 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | ND | 0.20 | 0.051 | ppbv | | ND | 0.87 | 0.22 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 0.22 | 0.20 | 0.051 | ppbv | | 0.96 | 0.87 | 0.22 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | ‡2 I | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 105% | | e | 55-128% |) | | | | | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



10 of 375

JC15931

4.1

Page 2 of 2

| Lab Sam Matrix: Method: Project: | AIR TO-1 | 931-2 - Soil Vapor 5 | r Comp. Summ sey Avenue, Bro | | 53 D P | Pate Sampled:0Pate Received:0Pate Received:< | |
|---|-------------|----------------------------|---------------------------------|-----|-----------|--|------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | 5W17103.D | 1 | 03/18/16 | TCH | n/a | n/a | V5W683 |
| | 5W17113.D | 1 | 03/18/16 | TCH | n/a | n/a | V5W684 |
| Run #2 | | | | | | | |

Report of Analysis

Run #2 20.0 ml

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|---------|------------------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 222 ^a | 4.0 | 0.72 | ppbv | 527 ^a | 9.5 | 1.7 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.11 | ppbv | ND | 1.8 | 0.24 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 1.1 | 0.80 | 0.13 | ppbv | 3.5 | 2.6 | 0.42 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.15 | ppbv | ND | 5.4 | 1.0 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.063 | ppbv | ND | 8.3 | 0.65 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.074 | ppbv | ND | 3.1 | 0.29 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.074 | ppbv | ND | 3.5 | 0.32 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.11 | ppbv | ND | 4.1 | 0.57 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.80 | 0.13 | ppbv | ND | 2.5 | 0.40 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.22 | ppbv | ND | 3.7 | 1.0 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.14 | ppbv | ND | 2.1 | 0.37 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.80 | 0.066 | ppbv | ND | 3.9 | 0.32 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.21 | ppbv | ND | 1.7 | 0.43 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.068 | ppbv | ND | 4.1 | 0.35 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.12 | ppbv | ND | 5.0 | 0.75 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | ND | 0.80 | 0.065 | ppbv | ND | 2.8 | 0.22 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.061 | ppbv | ND | 3.2 | 0.25 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.084 | ppbv | ND | 3.2 | 0.33 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.17 | ppbv | ND | 6.1 | 1.3 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.070 | 11 | ND | 3.2 | 0.28 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.088 | ppbv | ND | 3.7 | 0.41 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.18 | ppbv | ND | 2.9 | 0.65 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.56 | 0.80 | 0.076 | 11 | 2.8 | 4.0 | 0.38 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.21 | ppbv | ND | 6.8 | 1.8 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | ND | 3.2 | 0.44 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.084 | 11 | ND | 3.2 | 0.33 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.061 | ppbv | ND | 3.6 | 0.28 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.078 | * * | ND | 4.8 | 0.47 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.064 | * * | ND | 4.8 | 0.38 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | ND | 4.8 | 0.66 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.073 | ppbv | ND | 3.6 | 0.33 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 3



JC15931

E = Indicates value exceeds calibration range

| Report of Analysis | S | |
|--------------------|---|--|
|--------------------|---|--|

| Client Sample ID: | SV-10 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC15931-2 | Date Sampled: | 03/09/16 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A753 | Date Received: | 03/10/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| _ | | | |

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|---------|------|---------|-------|---|------------------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 46.9 | 2.0 | 0.30 | ppbv | | 88.4 | 3.8 | 0.57 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | ND | 0.80 | 0.17 | ppbv | | ND | 3.5 | 0.74 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 3.2 | 0.80 | 0.30 | ppbv | | 12 | 2.9 | 1.1 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 0.80 | 0.067 | ppbv | | ND | 3.9 | 0.33 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | | ppbv | | ND | 6.1 | 0.66 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.13 | ppbv | | ND | 5.6 | 0.91 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | ND | 0.80 | | ppbv | | ND | 3.3 | 0.33 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.081 | ppbv | | ND | 8.5 | 0.86 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 0.56 | 0.80 | | ppbv | J | 2.0 | 2.8 | 0.32 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 0.80 | 0.18 | ppbv | | ND | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 25.3 | 0.80 | 0.62 | ppbv | | 62.2 | 2.0 | 1.5 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | 0.43 | 0.80 | 0.10 | ppbv | J | 1.5 | 2.8 | 0.35 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 263 a | 4.0 | 0.97 | ppbv | | 776 ^a | 12 | 2.9 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | 1.2 | 0.80 | 0.22 | ppbv | | 4.9 | 3.3 | 0.90 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.079 | ppbv | | ND | 2.9 | 0.28 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.16 | ppbv | | ND | 3.3 | 0.66 | ug/m3 |
| 115-07-1 | 42 | Propylene | 2.6 | 2.0 | 0.13 | ppbv | | 4.5 | 3.4 | 0.22 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.062 | | | ND | 3.4 | 0.26 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.094 | ppbv | | ND | 4.4 | 0.51 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.064 | ppbv | | ND | 5.5 | 0.44 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.16 | ppbv | | ND | 4.4 | 0.87 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.22 | ppbv | | ND | 5.9 | 1.6 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | ND | 0.80 | 0.061 | ppbv | | ND | 3.9 | 0.30 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | 0.80 | 0.18 | ppbv | | ND | 3.9 | 0.88 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | ND | 0.80 | 0.091 | ppbv | | ND | 3.7 | 0.43 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 6.8 | 0.80 | 0.21 | ppbv | | 21 | 2.4 | 0.64 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 221 a | 0.80 | 0.46 | ppbv | | 1500 a | 5.4 | 3.1 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | 357 a | 4.0 | 0.91 | ppbv | | 1050 a | 12 | 2.7 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 1.6 | 0.80 | 0.050 | ppbv | | 6.0 | 3.0 | 0.19 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 0.80 | 0.16 | 0.074 | ppbv | | 4.3 | 0.86 | 0.40 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.088 | ppbv | | ND | 4.5 | 0.49 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.082 | ppbv | | ND | 2.0 | 0.21 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m, p-Xylene | 0.44 | 0.80 | 0.27 | ppbv | J | 1.9 | 3.5 | 1.2 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | ND | 0.80 | 0.20 | ppbv | | ND | 3.5 | 0.87 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 0.44 | 0.80 | 0.20 | ppbv | J | 1.9 | 3.5 | 0.87 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run | f 1 Rur | n# 2 | Limits | | | | | | |
| 460-00-4 | 4-Brom | nofluorobenzene 99% | 98% |) | 65-128% |) | | | | | |

ND = Not detectedMDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 3

12 of 375

JC15931

| | Report of Analysis | Page 3 of 3 |
|---|---|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-10Date Sampled:03/09/16JC15931-2Date Sampled:03/10/16AIR - Soil Vapor Comp.Summa ID:A753TO-15Date Received:03/10/162002-2024 Cropsey Avenue, Brooklyn, NYn/a | |
| VOA TO15 List CAS No. MW | Compound Result RL MDL Units Q Result RL | MDL Units |

(a) Result is from Run# 2

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



13 of 375 ACCUTEST JC15931

| Client Sa Lab Sam Matrix: Method: Project: | AII TO | 5931-3 R - Indoor Air -15 | r Comp. Summ sey Avenue, Bro | | 1 | Date Sampled: Date Received: Percent Solids: | |
|--|-------------------------------|---------------------------------|---------------------------------|------------------|-------------------------|--|------------------------------|
| Run #1 Run #2 | File ID 5W17130.D | DF 1 | Analyzed 03/19/16 | By TCH | Prep Date n/a | Prep Batc n/a | h Analytical Batch V5W684 |
| | Initial Volu 400 ml | me | | | | | |

Report of Analysis

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|------|-------|---------|--------|------|-------|-------|
| 67-64-1 | 58.08 | Acetone | 6.6 | 0.20 | 0.036 | ppbv | 16 | 0.48 | 0.086 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.20 | 0.028 | ppbv | ND | 0.44 | 0.062 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 0.80 | 0.20 | 0.031 | ppbv | 2.6 | 0.64 | 0.099 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.20 | 0.039 | | ND | 1.3 | 0.26 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.20 | 0.016 | | ND | 2.1 | 0.17 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.20 | 0.018 | ppbv | ND | 0.78 | 0.070 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.20 | 0.018 | ppbv | ND | 0.87 | 0.079 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.20 | 0.027 | ppbv | ND | 1.0 | 0.14 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.20 | 0.031 | ppbv | ND | 0.62 | 0.097 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.20 | 0.056 | ppbv | ND | 0.92 | 0.26 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.20 | 0.036 | | ND | 0.53 | 0.095 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.20 | 0.017 | | ND | 0.98 | 0.083 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | 0.87 | 0.20 | 0.052 | ppbv | 1.8 | 0.41 | 0.11 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.20 | 0.027 | ppbv | ND | 0.63 | 0.085 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.20 | 0.017 | ppbv | ND | 1.0 | 0.088 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.20 | 0.031 | | ND | 1.3 | 0.20 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | ND | 0.20 | 0.016 | | ND | 0.69 | 0.055 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.20 | 0.015 | | ND | 0.81 | 0.061 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.20 | 0.021 | | ND | 0.79 | 0.083 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.20 | 0.042 | ppbv | ND | 1.5 | 0.32 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.20 | 0.018 | ppbv | ND | 0.81 | 0.073 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.20 | 0.022 | ppbv | ND | 0.92 | 0.10 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.20 | 0.045 | ppbv | ND | 0.72 | 0.16 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.64 | 0.20 | 0.019 | ppbv | 3.2 | 0.99 | 0.094 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.20 | 0.053 | | ND | 1.7 | 0.45 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.20 | 0.028 | | ND | 0.79 | 0.11 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.20 | | | ND | 0.79 | 0.083 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.20 | 0.015 | | ND | 0.91 | 0.068 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.20 | 0.020 | | ND | 1.2 | 0.12 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.20 | 0.016 | | ND | 1.2 | 0.096 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.20 | 0.027 | 11 | ND | 1.2 | 0.16 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.20 | 0.018 | ppbv | ND | 0.91 | 0.082 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 1 of 2

14 of 375

ACCUTEST JC15931

SGS

| Report or marysis | Report | of | Anal | lysis |
|-------------------|--------|----|------|-------|
|-------------------|--------|----|------|-------|

| Client Sample ID: | IAQ-10 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC15931-3 | Date Sampled: | 03/09/16 |
| Matrix: | AIR - Indoor Air Comp. Summa ID: A371 | Date Received: | 03/10/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | · · · | | |

VOA TO15 List

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| 100-41-4 10 141-78-6 88 622-96-8 12 76-13-1 18 76-14-2 17 142-82-5 10 87-68-3 26 110-54-3 86 591-78-6 10 67-63-0 60 | 8 20.2 87.4 70.9 00.2 60.8 66.17 00 | Ethanol Ethylbenzene Ethyl Acetate 4-Ethyltoluene Freon 113 Freon 114 Heptane Hexachlorobutadiene Hexane | 18.6 0.13 1.9 ND ND ND 0.19 ND | 0.50 0.20 0.20 0.20 0.20 0.20 0.20 0.20 | 0.075 0.042 0.075 0.017 0.021 | ppbv ppbv ppbv | J | 35.0 0.56 6.8 ND | 0.94 0.87 0.72 0.98 | 0.14 0.18 0.27 0.084 | ug/m3 ug/m3 ug/m3 |
|---|--|--|---|--|---|----------------------|---|---------------------------|------------------------------|-------------------------------|-------------------------|
| 141-78-6 88 622-96-8 12 76-13-1 18 76-14-2 17 142-82-5 10 87-68-3 26 110-54-3 86 591-78-6 10 67-63-0 60 | 8 20.2 87.4 70.9 00.2 60.8 66.17 00 | Ethyl Acetate 4-Ethyltoluene Freon 113 Freon 114 Heptane Hexachlorobutadiene Hexane | 1.9 ND ND ND 0.19 | 0.20 0.20 0.20 0.20 | 0.075 0.017 0.021 | ppbv ppbv | J | 6.8 ND | 0.72 | 0.27 | |
| 622-96-8 12 76-13-1 18 76-14-2 17 142-82-5 10 87-68-3 26 110-54-3 86 591-78-6 10 67-63-0 60 | 20.2 87.4 70.9 00.2 60.8 66.17 00 | 4-Ethyltoluene Freon 113 Freon 114 Heptane Hexachlorobutadiene Hexane | ND ND ND 0.19 | 0.20 0.20 0.20 | 0.017 0.021 | ppbv | | ND | | | ug/m3 |
| 76-13-1 18 76-14-2 17 142-82-5 10 87-68-3 26 110-54-3 86 591-78-6 10 67-63-0 60 | 87.4 70.9 00.2 60.8 6.17 00 | Freon 113 Freon 114 Heptane Hexachlorobutadiene Hexane | ND ND 0.19 | 0.20 0.20 | 0.021 | | | | 0.98 | 0 001 | |
| 76-14-2 17 142-82-5 10 87-68-3 26 110-54-3 86 591-78-6 10 67-63-0 60 | 70.9 00.2 60.8 6.17 00 | Freon 114 Heptane Hexachlorobutadiene Hexane | ND 0.19 | 0.20 | | ppbv | | ND | | 0.064 | ug/m3 |
| 142-82-5 10 87-68-3 26 110-54-3 86 591-78-6 10 67-63-0 60 | 00.2 60.8 6.17 00 | Heptane Hexachlorobutadiene Hexane | 0.19 | | 0.021 | | | ND | 1.5 | 0.16 | ug/m3 |
| 87-68-3 26 110-54-3 86 591-78-6 10 67-63-0 60 | 260.8 6.17 00 | Hexachlorobutadiene Hexane | | 0.20 | 0.051 | ppbv | | ND | 1.4 | 0.22 | ug/m3 |
| 110-54-386591-78-61067-63-060 | 6.17 00 | Hexane | ND | | 0.020 | | J | 0.78 | 0.82 | 0.082 | ug/m3 |
| 591-78-6 10 67-63-0 60 | 00 | | | 0.20 | | ppbv | | ND | 2.1 | 0.21 | ug/m3 |
| 67-63-0 60 | | | 0.69 | 0.20 | 0.023 | | | 2.4 | 0.70 | 0.081 | ug/m3 |
| | 0.1 | 2-Hexanone | ND | 0.20 | | ppbv | | ND | 0.82 | 0.18 | ug/m3 |
| 75.00.2 | | Isopropyl Alcohol | 1.8 | 0.20 | 0.16 | ppbv | | 4.4 | 0.49 | 0.39 | ug/m3 |
| | | Methylene chloride | 1.6 | 0.20 | 0.025 | ppbv | | 5.6 | 0.69 | 0.087 | ug/m3 |
| 78-93-3 72 | | Methyl ethyl ketone | 0.75 | 0.20 | | ppbv | | 2.2 | 0.59 | 0.14 | ug/m3 |
| | | Methyl Isobutyl Ketone | ND | 0.20 | 0.055 | | | ND | 0.82 | 0.23 | ug/m3 |
| | | Methyl Tert Butyl Ether | ND | 0.20 | 0.020 | | | ND | 0.72 | | ug/m3 |
| | | Methylmethacrylate | ND | 0.20 | 0.040 | | | ND | 0.82 | 0.16 | ug/m3 |
| 115-07-1 42 | | Propylene | 3.2 | 0.50 | 0.032 | | | 5.5 | 0.86 | | ug/m3 |
| | | Styrene | ND | 0.20 | 0.015 | | | ND | 0.85 | 0.064 | ug/m3 |
| 71-55-6 13 | 33.4 | 1,1,1-Trichloroethane | ND | 0.20 | 0.024 | | | ND | 1.1 | 0.13 | ug/m3 |
| | 67.9 | 1,1,2,2-Tetrachloroethane | ND | 0.20 | 0.016 | | | ND | 1.4 | 0.11 | ug/m3 |
| | 33.4 | 1,1,2-Trichloroethane | ND | 0.20 | 0.039 | | | ND | 1.1 | 0.21 | ug/m3 |
| | | 1,2,4-Trichlorobenzene | ND | 0.20 | 0.056 | | | ND | 1.5 | 0.42 | ug/m3 |
| | | 1,2,4-Trimethylbenzene | ND | 0.20 | 0.015 | | | ND | 0.98 | 0.074 | ug/m3 |
| | 20.2 | 1,3,5-Trimethylbenzene | ND | 0.20 | 0.045 | | | ND | 0.98 | 0.22 | ug/m3 |
| | | 2,2,4-Trimethylpentane | 0.49 | 0.20 | 0.023 | | | 2.3 | 0.93 | 0.11 | ug/m3 |
| | | Tertiary Butyl Alcohol | ND | 0.20 | 0.053 | | | ND | 0.61 | 0.16 | ug/m3 |
| | | Tetrachloroethylene | 0.11 | | 0.023 | | | 0.75 | 0.27 | 0.16 | ug/m3 |
| | | Tetrahydrofuran | 0.52 | 0.20 | 0.045 | | | 1.5 | 0.59 | 0.13 | ug/m3 |
| | | Toluene | 1.3 | 0.20 | 0.012 | | | 4.9 | 0.75 | 0.045 | ug/m3 |
| | | Trichloroethylene | ND | | 0.019 | ~ ~ | | ND | 0.21 | 0.10 | ug/m3 |
| | | Trichlorofluoromethane | 0.34 | 0.20 | 0.022 | | | 1.9 | 1.1 | 0.12 | ug/m3 |
| | 2.5 | Vinyl chloride | ND | 0.20 | 0.021 | | | ND | 0.51 | 0.054 | ug/m3 |
| 108-05-4 86 | | Vinyl Acetate | ND | 0.20 | 0.054 | | | ND | 0.70 | 0.19 | ug/m3 |
| | | m,p-Xylene | 0.40 | 0.20 | 0.068 | | | 1.7 | 0.87 | 0.30 | ug/m3 |
| | | o-Xylene | 0.15 | 0.20 | 0.051 | | J | 0.65 | 0.87 | 0.22 | ug/m3 |
| 1330-20-7 10 | 06.2 | Xylenes (total) | 0.54 | 0.20 | 0.051 | ppbv | | 2.3 | 0.87 | 0.22 | ug/m3 |
| CAS No. Su | burroga | ate Recoveries Run# | 1 Run# | 2 L | imits | | | | | | |
| 460-00-4 4- | -Brom | ofluorobenzene 98% | | 6 | 5-128% | | | | | | |

ND = Not detectedMDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Page 2 of 2

4.3 4



15 of 375

JC15931



Section 4

Sample Results

Report of Analysis





| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC165 AIR - TO-15 | 571-1 Soil Vapor 5 | Comp. Summ ey Avenue, Bro | | | Date Sampled:03/17/16Date Received:03/18/16Percent Solids:n/a | | | | |
|---|---|--------------------------|------------------------------|-----|-----------|---|--------------------|--|--|--|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batc | h Analytical Batch | | | |
| Run #1 | 3W52959.D | 1 | 03/24/16 | YMH | n/a | n/a | V3W1999 | | | |
| Run #2 | 3W52993.D | 1.38 | 03/26/16 | YMH | n/a | n/a | V3W2000 | | | |
| Run #1 Run #2 | Initial Volum 100 ml 20.0 ml | e | | | | | | | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|---------|-------------------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 551 ^a | 5.5 | 1.0 | ppbv | 1310 ^a | 13 | 2.4 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.11 | ppbv | ND | 1.8 | 0.24 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 4.9 | 0.80 | 0.13 | ppbv | 16 | 2.6 | 0.42 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.15 | ppbv | ND | 5.4 | 1.0 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.063 | ppbv | ND | 8.3 | 0.65 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.074 | ppbv | ND | 3.1 | 0.29 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.074 | ppbv | ND | 3.5 | 0.32 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.11 | ppbv | ND | 4.1 | 0.57 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | 1.9 | 0.80 | 0.13 | ppbv | 5.9 | 2.5 | 0.40 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.22 | ppbv | ND | 3.7 | 1.0 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.14 | ppbv | ND | 2.1 | 0.37 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | 3.5 | 0.80 | 0.066 | ppbv | 17 | 3.9 | 0.32 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.21 | ppbv | ND | 1.7 | 0.43 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.068 | ppbv | ND | 4.1 | 0.35 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.12 | ppbv | ND | 5.0 | 0.75 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 2.3 | 0.80 | 0.065 | ppbv | 7.9 | 2.8 | 0.22 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.061 | ppbv | ND | 3.2 | 0.25 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.084 | ppbv | ND | 3.2 | 0.33 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.17 | ppbv | ND | 6.1 | 1.3 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.070 | ppbv | ND | 3.2 | 0.28 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.088 | ppbv | ND | 3.7 | 0.41 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.18 | ppbv | ND | 2.9 | 0.65 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.59 | 0.80 | 0.076 | ppbv J | 2.9 | 4.0 | 0.38 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.21 | ppbv | ND | 6.8 | 1.8 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | ND | 3.2 | 0.44 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.084 | ppbv | ND | 3.2 | 0.33 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.061 | ppbv | ND | 3.6 | 0.28 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.078 | ppbv | ND | 4.8 | 0.47 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.064 | ppbv | ND | 4.8 | 0.38 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | ND | 4.8 | 0.66 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.073 | ppbv | ND | 3.6 | 0.33 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 3

11 of 658

ACCUTEST JC16571

E = Indicates value exceeds calibration range

| Report of A | nalysis |
|-------------|---------|
|-------------|---------|

| Client Sample ID: | SV-14 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC16571-1 | Date Sampled: | 03/17/16 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A752 | Date Received: | 03/18/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| _ | | | |

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|------|---------|---------|--------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 19.7 | 2.0 | 0.30 | ppbv | 37.1 | 3.8 | 0.57 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 16.9 | 0.80 | 0.17 | ppbv | 73.4 | 3.5 | 0.74 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | ND | 0.80 | 0.30 | ppbv | ND | 2.9 | 1.1 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | 3.8 | 0.80 | 0.067 | ppbv | 19 | 3.9 | 0.33 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.086 | ppbv | ND | 6.1 | 0.66 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.13 | ppbv | ND | 5.6 | 0.91 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 15.1 | 0.80 | 0.081 | ppbv | 61.9 | 3.3 | 0.33 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.081 | ppbv | ND | 8.5 | 0.86 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 6.3 | 0.80 | 0.090 | ppbv | 22 | 2.8 | 0.32 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | 2.2 | 0.80 | 0.18 | ppbv | 9.0 | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 3.8 | 0.80 | 0.62 | ppbv | 9.3 | 2.0 | 1.5 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.10 | ppbv | ND | 2.8 | 0.35 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 12.3 | 0.80 | 0.19 | ppbv | 36.3 | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | 1.2 | 0.80 | 0.22 | ppbv | 4.9 | 3.3 | 0.90 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.079 | ppbv | ND | 2.9 | 0.28 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.16 | ppbv | ND | 3.3 | 0.66 | ug/m3 |
| 115-07-1 | 42 | Propylene | 6.9 | 2.0 | 0.13 | ppbv | 12 | 3.4 | 0.22 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.062 | | ND | 3.4 | 0.26 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.094 | ppbv | ND | 4.4 | 0.51 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.064 | ppbv | ND | 5.5 | 0.44 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.16 | ppbv | ND | 4.4 | 0.87 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.22 | ppbv | ND | 5.9 | 1.6 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 12.8 | 0.80 | 0.061 | ppbv | 62.9 | 3.9 | 0.30 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 3.8 | 0.80 | 0.18 | ppbv | 19 | 3.9 | 0.88 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 2.4 | 0.80 | 0.091 | ppbv | 11 | 3.7 | 0.43 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 2.5 | 0.80 | 0.21 | ppbv | 7.6 | 2.4 | 0.64 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 38.0 | 0.16 | 0.092 | ppbv | 258 | 1.1 | 0.62 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.80 | 0.18 | ppbv | ND | 2.4 | 0.53 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 91.2 | 0.80 | 0.050 | ppbv | 344 | 3.0 | 0.19 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | 1.4 | 0.16 | 0.074 | ppbv | 7.5 | 0.86 | 0.40 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.088 | ppbv | ND | 4.5 | 0.49 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.082 | ppbv | ND | 2.0 | 0.21 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 60.8 | 0.80 | 0.27 | ppbv | 264 | 3.5 | 1.2 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 17.6 | 0.80 | 0.20 | ppbv | 76.4 | 3.5 | 0.87 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 78.3 | 0.80 | 0.20 | ppbv | 340 | 3.5 | 0.87 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run | # 2 | Limits | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 91% | 93% | | 65-128% |) | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound

12 of 658 SGS ACCUTEST

JC16571

Page 2 of 3

4.1 **4**

| | Report of Analy | vsis | Page 3 of 3 |
|---|--|---|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-14 JC16571-1 AIR - Soil Vapor Comp. Summa ID: A752 TO-15 2002-2024 Cropsey Avenue, Brooklyn, NY | Date Sampled:03/17/16Date Received:03/18/16Percent Solids:n/a | |
| VOA TO15 List | | | |
| CAS No. MW | Compound Result RL | MDL Units Q Result RL | MDL Units |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



13 of 658

JC16571

| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC165 AIR - TO-15 | 571-2 Soil Vapor 5 | Comp. Summ | | | Date Sampled:03/17/16Date Received:03/18/16Percent Solids:n/a | | | | |
|---|---|--------------------------|------------|-----|-----------|---|--------------------|--|--|--|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batcl | h Analytical Batch | | | |
| Run #1 | 3W52960.D | 1 | 03/24/16 | YMH | n/a | n/a | V3W1999 | | | |
| Run #2 | 3W52994.D | 1 | 03/26/16 | YMH | n/a | n/a | V3W2000 | | | |
| Run #1 Run #2 | Initial Volum 100 ml 50.0 ml | е | | | | | | | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|------------------|------|-------|---------|------------------|-----|------|-------|
| 67-64-1 | 58.08 | Acetone | 188 ^a | 1.6 | 0.29 | ppbv | 447 ^a | 3.8 | 0.69 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.80 | 0.11 | ppbv | ND | 1.8 | 0.24 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 4.6 | 0.80 | 0.13 | ppbv | 15 | 2.6 | 0.42 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.80 | 0.15 | ppbv | ND | 5.4 | 1.0 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.80 | 0.063 | ppbv | ND | 8.3 | 0.65 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.80 | 0.074 | ppbv | ND | 3.1 | 0.29 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.80 | 0.074 | ppbv | ND | 3.5 | 0.32 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.80 | 0.11 | ppbv | ND | 4.1 | 0.57 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | 2.0 | 0.80 | 0.13 | ppbv | 6.2 | 2.5 | 0.40 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.80 | 0.22 | ppbv | ND | 3.7 | 1.0 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.80 | 0.14 | ppbv | ND | 2.1 | 0.37 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | 11.8 | 0.80 | 0.066 | ppbv | 57.6 | 3.9 | 0.32 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 0.80 | 0.21 | ppbv | ND | 1.7 | 0.43 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.80 | 0.11 | ppbv | ND | 2.5 | 0.34 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.80 | 0.068 | ppbv | ND | 4.1 | 0.35 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.80 | 0.12 | ppbv | ND | 5.0 | 0.75 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 2.9 | 0.80 | 0.065 | ppbv | 10 | 2.8 | 0.22 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.80 | 0.061 | ppbv | ND | 3.2 | 0.25 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.80 | 0.084 | ppbv | ND | 3.2 | 0.33 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.80 | 0.17 | ppbv | ND | 6.1 | 1.3 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.80 | 0.070 | ppbv | ND | 3.2 | 0.28 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.80 | 0.088 | ppbv | ND | 3.7 | 0.41 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.80 | 0.18 | ppbv | ND | 2.9 | 0.65 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.58 | 0.80 | 0.076 | ppbv J | 2.9 | 4.0 | 0.38 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.80 | 0.21 | ppbv | ND | 6.8 | 1.8 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.80 | 0.11 | ppbv | ND | 3.2 | 0.44 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.80 | 0.084 | ppbv | ND | 3.2 | 0.33 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.80 | 0.061 | ppbv | ND | 3.6 | 0.28 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.80 | 0.078 | ppbv | ND | 4.8 | 0.47 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.80 | 0.064 | ppbv | ND | 4.8 | 0.38 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.80 | 0.11 | ppbv | ND | 4.8 | 0.66 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.80 | 0.073 | ppbv | ND | 3.6 | 0.33 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 3

14 of 658

ACCUTEST JC16571

E = Indicates value exceeds calibration range

| Report of A | nalysis |
|-------------|---------|
|-------------|---------|

| Client Sample ID: | SV-13 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC16571-2 | Date Sampled: | 03/17/16 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A219 | Date Received: | 03/18/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|------------|---------|-------|---|--------|------|------|-------|
| 64-17-5 | 46.07 | Ethanol | 17.9 | 2.0 | 0.30 | ppbv | | 33.7 | 3.8 | 0.57 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 18.3 | 0.80 | 0.17 | ppbv | | 79.5 | 3.5 | 0.74 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | ND | 0.80 | 0.30 | ppbv | | ND | 2.9 | 1.1 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | 4.7 | 0.80 | 0.067 | ppbv | | 23 | 3.9 | 0.33 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 0.80 | 0.086 | ppbv | | ND | 6.1 | 0.66 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 0.80 | 0.13 | ppbv | | ND | 5.6 | 0.91 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 15.6 | 0.80 | 0.081 | ppbv | | 63.9 | 3.3 | 0.33 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 0.80 | 0.081 | ppbv | | ND | 8.5 | 0.86 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 5.6 | 0.80 | 0.090 | ppbv | | 20 | 2.8 | 0.32 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | 0.61 | 0.80 | 0.18 | ppbv | J | 2.5 | 3.3 | 0.74 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 2.1 | 0.80 | 0.62 | ppbv | | 5.2 | 2.0 | 1.5 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 0.80 | 0.10 | ppbv | | ND | 2.8 | 0.35 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | 7.7 | 0.80 | 0.19 | ppbv | | 23 | 2.4 | 0.56 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | 2.1 | 0.80 | 0.22 | ppbv | | 8.6 | 3.3 | 0.90 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 0.80 | 0.079 | ppbv | | ND | 2.9 | 0.28 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 0.80 | 0.16 | ppbv | | ND | 3.3 | 0.66 | ug/m3 |
| 115-07-1 | 42 | Propylene | 3.2 | 2.0 | 0.13 | ppbv | | 5.5 | 3.4 | 0.22 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 0.80 | 0.062 | ppbv | | ND | 3.4 | 0.26 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 0.80 | 0.094 | ppbv | | ND | 4.4 | 0.51 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 0.80 | 0.064 | ppbv | | ND | 5.5 | 0.44 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 0.80 | 0.16 | ppbv | | ND | 4.4 | 0.87 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 0.80 | 0.22 | ppbv | | ND | 5.9 | 1.6 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | 16.3 | 0.80 | 0.061 | ppbv | | 80.1 | 3.9 | 0.30 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | 4.8 | 0.80 | 0.18 | ppbv | | 24 | 3.9 | 0.88 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | 2.8 | 0.80 | 0.091 | ppbv | | 13 | 3.7 | 0.43 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | 1.2 | 0.80 | 0.21 | ppbv | | 3.6 | 2.4 | 0.64 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | 2.1 | 0.16 | 0.092 | ppbv | | 14 | 1.1 | 0.62 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 0.80 | 0.18 | ppbv | | ND | 2.4 | 0.53 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 94.8 | 0.80 | 0.050 | ppbv | | 357 | 3.0 | 0.19 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | 0.16 | 0.074 | ppbv | | ND | 0.86 | 0.40 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 0.80 | 0.088 | ppbv | | ND | 4.5 | 0.49 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | 0.80 | 0.082 | ppbv | | ND | 2.0 | 0.21 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 0.80 | 0.22 | ppbv | | ND | 2.8 | 0.77 | ug/m3 |
| | 106.2 | m,p-Xylene | 68.4 | 0.80 | 0.27 | ppbv | | 297 | 3.5 | 1.2 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 20.1 | 0.80 | 0.20 | ppbv | | 87.3 | 3.5 | 0.87 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 88.5 | 0.80 | 0.20 | ppbv | | 384 | 3.5 | 0.87 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run# | # 2 | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 92% | 93% | | 65-128% |) | | | | | |

ND = Not detectedMDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





15 of 658

JC16571

| | Report of Analysis | Page 3 of 3 |
|---|---|-------------|
| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | SV-13 JC16571-2 Date Sampled: 03/17/16 AIR - Soil Vapor Comp. Summa ID: A219 Date Received: 03/18/16 TO-15 Percent Solids: n/a 2002-2024 Cropsey Avenue, Brooklyn, NY | 4.2 |
| VOA TO15 List | | |
| CAS No. MW | Compound Result RL MDL Units Q Result RL | MDL Units |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



16 of 658 ACCUTEST JC16571

| | | | Repo | ort of A | nalysis | | Page 1 of 2 |
|--|------------------------------|---------------------|---------------------------------|------------------|-------------------------|--------------------------|-----------------------------|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC Al |)-15 | r Comp. Sumn sey Avenue, Bro | | | L | 03/17/16 03/18/16 n/a |
| Run #1 Run #2 | File ID 3W52961. | DF D 59.2 | Analyzed 03/24/16 | By YMH | Prep Date n/a | Prep Batch n/a | Analytical Batch V3W1999 |
| Run #1 | Initial Vol 200 ml | ume | | | | | |

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|----|-----|-------|---|--------|-----|-----|-------|
| 67-64-1 | 58.08 | Acetone | ND | 24 | 4.3 | ppbv | | ND | 57 | 10 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 24 | 3.3 | ppbv | | ND | 53 | 7.3 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | 10.6 | 24 | 3.7 | ppbv | J | 33.9 | 77 | 12 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 24 | 4.6 | ppbv | | ND | 160 | 31 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 24 | 1.9 | ppbv | | ND | 250 | 20 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 24 | 2.2 | ppbv | | ND | 93 | 8.5 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 24 | 2.2 | ppbv | | ND | 100 | 9.6 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 24 | 3.2 | ppbv | | ND | 120 | 16 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 24 | 3.7 | ppbv | | ND | 75 | 12 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 24 | 6.6 | ppbv | | ND | 110 | 30 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 24 | 4.2 | ppbv | | ND | 63 | 11 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 24 | 2.0 | ppbv | | ND | 120 | 9.8 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | ND | 24 | 6.2 | ppbv | | ND | 50 | 13 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 24 | 3.1 | ppbv | | ND | 75 | 9.7 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 24 | 2.0 | ppbv | | ND | 120 | 10 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 24 | 3.7 | ppbv | | ND | 150 | 23 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | 115 | 24 | 1.9 | ppbv | | 396 | 83 | 6.5 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 24 | 1.8 | ppbv | | ND | 97 | 7.3 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 24 | 2.5 | ppbv | | ND | 95 | 9.9 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 24 | 5.0 | ppbv | | ND | 180 | 38 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 24 | 2.1 | ppbv | | ND | 97 | 8.5 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 24 | 2.6 | ppbv | | ND | 110 | 12 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 24 | 5.3 | ppbv | | ND | 86 | 19 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | ND | 24 | 2.3 | ppbv | | ND | 120 | 11 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 24 | 6.2 | ppbv | | ND | 200 | 53 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 24 | 3.3 | ppbv | | ND | 95 | 13 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | 30.2 | 24 | 2.5 | ppbv | | 120 | 95 | 9.9 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 24 | 1.8 | ppbv | | ND | 110 | 8.2 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 24 | 2.3 | ppbv | | ND | 140 | 14 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 24 | 1.9 | ppbv | | ND | 140 | 11 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 24 | 3.2 | ppbv | | ND | 140 | 19 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 24 | 2.2 | ppbv | | ND | 110 | 10 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



4.3



E = Indicates value exceeds calibration range

| Client Sample ID: | SV-12 | | |
|-------------------|---|-----------------|----------|
| Lab Sample ID: | JC16571-3 | Date Sampled: | 03/17/16 |
| Matrix: | AIR - Soil Vapor Comp. Summa ID: A364, M206 | Date Received: | 03/18/16 |
| Method: | TO-15 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|--------|-----|---------|-------|---|--------|-----|-----|-------|
| 64-17-5 | 46.07 | Ethanol | ND | 59 | 8.9 | ppbv | | ND | 110 | 17 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | 11.6 | 24 | 5.0 | ppbv | J | 50.4 | 100 | 22 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | ND | 24 | 8.8 | ppbv | | ND | 86 | 32 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | 24 | 2.0 | ppbv | | ND | 120 | 9.8 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | 24 | 2.5 | ppbv | | ND | 180 | 19 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | 24 | 3.7 | ppbv | | ND | 170 | 26 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | 24.7 | 24 | 2.4 | ppbv | | 101 | 98 | 9.8 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | 24 | 2.4 | ppbv | | ND | 260 | 26 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 488 | 24 | 2.7 | ppbv | | 1720 | 85 | 9.5 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | 24 | 5.3 | ppbv | | ND | 98 | 22 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | ND | 24 | 18 | ppbv | | ND | 59 | 44 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | ND | 24 | 2.9 | ppbv | | ND | 83 | 10 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | ND | 24 | 5.7 | ppbv | | ND | 71 | 17 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | 24 | 6.5 | ppbv | | ND | 98 | 27 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | 24 | 2.3 | ppbv | | ND | 87 | 8.3 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | 24 | 4.7 | ppbv | | ND | 98 | 19 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | 59 | 3.8 | ppbv | | ND | 100 | 6.5 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | 24 | 1.8 | ppbv | | ND | 100 | 7.7 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | 24 | 2.8 | ppbv | | ND | 130 | 15 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | 24 | 1.9 | ppbv | | ND | 160 | 13 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | 24 | 4.6 | ppbv | | ND | 130 | 25 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | 24 | 6.6 | ppbv | | ND | 180 | 49 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | ND | 24 | 1.8 | ppbv | | ND | 120 | 8.8 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | 24 | 5.3 | ppbv | | ND | 120 | 26 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | ND | 24 | 2.7 | ppbv | | ND | 110 | 13 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | ND | 24 | 6.3 | ppbv | | ND | 73 | 19 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | ND | 4.7 | 2.7 | ppbv | | ND | 32 | 18 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | 24 | 5.4 | ppbv | | ND | 71 | 16 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 52.8 | 24 | 1.5 | ppbv | | 199 | 90 | 5.7 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | 4.7 | 2.2 | ppbv | | ND | 25 | 12 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | 24 | 2.6 | ppbv | | ND | 130 | 15 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | 334 | 24 | 2.4 | ppbv | | 854 | 61 | 6.1 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | 24 | 6.4 | ppbv | | ND | 84 | 23 | ug/m3 |
| | 106.2 | m,p-Xylene | 41.8 | 24 | 8.1 | ppbv | | 182 | 100 | 35 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | 12.5 | 24 | 6.0 | ppbv | J | 54.3 | 100 | 26 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | 54.3 | 24 | 6.0 | ppbv | | 236 | 100 | 26 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 Run | # 2 | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 88% | | | 65-128% |) | | | | | |

ND = Not detectedMDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 2

4.3

4



JC16571

SGS

| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC165 AIR - TO-15 | Ambient A | | nma ID: M ooklyn, NY | | Date Sampled: Date Received: Percent Solids: | |
|---|---------------------------------|----------------|--------------------------|-------------------------|-------------------------|--|-------------------------------|
| Run #1 Run #2 | File ID 3W52962.D | DF 1 | Analyzed 03/25/16 | By YMH | Prep Date n/a | Prep Batc n/a | h Analytical Batch V3W1999 |
| Run #1 | Initial Volume 400 ml | 2 | | | | | |

Report of Analysis

Run #2

VOA TO15 List

| CAS No. | MW | Compound | Result | RL | MDL | Units | Q | Result | RL | MDL | Units |
|------------|-------|----------------------------|--------|------|-------|-------|---|--------|------|-------|-------|
| 67-64-1 | 58.08 | Acetone | 0.85 | 0.20 | 0.036 | ppbv | | 2.0 | 0.48 | 0.086 | ug/m3 |
| 106-99-0 | 54.09 | 1,3-Butadiene | ND | 0.20 | 0.028 | | | ND | 0.44 | 0.062 | ug/m3 |
| 71-43-2 | 78.11 | Benzene | ND | 0.20 | 0.031 | ppbv | | ND | 0.64 | 0.099 | ug/m3 |
| 75-27-4 | 163.8 | Bromodichloromethane | ND | 0.20 | 0.039 | ppbv | | ND | 1.3 | 0.26 | ug/m3 |
| 75-25-2 | 252.8 | Bromoform | ND | 0.20 | 0.016 | | | ND | 2.1 | 0.17 | ug/m3 |
| 74-83-9 | 94.94 | Bromomethane | ND | 0.20 | 0.018 | ppbv | | ND | 0.78 | 0.070 | ug/m3 |
| 593-60-2 | 106.9 | Bromoethene | ND | 0.20 | 0.018 | ppbv | | ND | 0.87 | 0.079 | ug/m3 |
| 100-44-7 | 126 | Benzyl Chloride | ND | 0.20 | 0.027 | ppbv | | ND | 1.0 | 0.14 | ug/m3 |
| 75-15-0 | 76.14 | Carbon disulfide | ND | 0.20 | 0.031 | ppbv | | ND | 0.62 | 0.097 | ug/m3 |
| 108-90-7 | 112.6 | Chlorobenzene | ND | 0.20 | 0.056 | ppbv | | ND | 0.92 | 0.26 | ug/m3 |
| 75-00-3 | 64.52 | Chloroethane | ND | 0.20 | 0.036 | | | ND | 0.53 | 0.095 | ug/m3 |
| 67-66-3 | 119.4 | Chloroform | ND | 0.20 | 0.017 | | | ND | 0.98 | 0.083 | ug/m3 |
| 74-87-3 | 50.49 | Chloromethane | 0.19 | 0.20 | 0.052 | ~ ~ | J | 0.39 | 0.41 | 0.11 | ug/m3 |
| 107-05-1 | 76.53 | 3-Chloropropene | ND | 0.20 | 0.027 | | | ND | 0.63 | 0.085 | ug/m3 |
| 95-49-8 | 126.6 | 2-Chlorotoluene | ND | 0.20 | 0.017 | | | ND | 1.0 | 0.088 | ug/m3 |
| 56-23-5 | 153.8 | Carbon tetrachloride | ND | 0.20 | 0.031 | | | ND | 1.3 | 0.20 | ug/m3 |
| 110-82-7 | 84.16 | Cyclohexane | ND | 0.20 | 0.016 | | | ND | 0.69 | 0.055 | ug/m3 |
| 75-34-3 | 98.96 | 1,1-Dichloroethane | ND | 0.20 | 0.015 | | | ND | 0.81 | 0.061 | ug/m3 |
| 75-35-4 | 96.94 | 1,1-Dichloroethylene | ND | 0.20 | 0.021 | | | ND | 0.79 | 0.083 | ug/m3 |
| 106-93-4 | 187.9 | 1,2-Dibromoethane | ND | 0.20 | 0.042 | | | ND | 1.5 | 0.32 | ug/m3 |
| 107-06-2 | 98.96 | 1,2-Dichloroethane | ND | 0.20 | 0.018 | | | ND | 0.81 | 0.073 | ug/m3 |
| 78-87-5 | 113 | 1,2-Dichloropropane | ND | 0.20 | 0.022 | | | ND | 0.92 | 0.10 | ug/m3 |
| 123-91-1 | 88.12 | 1,4-Dioxane | ND | 0.20 | 0.045 | ppbv | | ND | 0.72 | 0.16 | ug/m3 |
| 75-71-8 | 120.9 | Dichlorodifluoromethane | 0.14 | 0.20 | 0.019 | ppbv | J | 0.69 | 0.99 | 0.094 | ug/m3 |
| 124-48-1 | 208.3 | Dibromochloromethane | ND | 0.20 | 0.053 | ppbv | | ND | 1.7 | 0.45 | ug/m3 |
| 156-60-5 | 96.94 | trans-1,2-Dichloroethylene | ND | 0.20 | 0.028 | | | ND | 0.79 | 0.11 | ug/m3 |
| 156-59-2 | 96.94 | cis-1,2-Dichloroethylene | ND | 0.20 | 0.021 | | | ND | 0.79 | 0.083 | ug/m3 |
| 10061-01-5 | 111 | cis-1,3-Dichloropropene | ND | 0.20 | 0.015 | | | ND | 0.91 | 0.068 | ug/m3 |
| 541-73-1 | 147 | m-Dichlorobenzene | ND | 0.20 | 0.020 | | | ND | 1.2 | 0.12 | ug/m3 |
| 95-50-1 | 147 | o-Dichlorobenzene | ND | 0.20 | 0.016 | | | ND | 1.2 | 0.096 | ug/m3 |
| 106-46-7 | 147 | p-Dichlorobenzene | ND | 0.20 | 0.027 | 11 | | ND | 1.2 | 0.16 | ug/m3 |
| 10061-02-6 | 111 | trans-1,3-Dichloropropene | ND | 0.20 | 0.018 | ppbv | | ND | 0.91 | 0.082 | ug/m3 |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





| Client Sample ID: | AMBIENT UPWIND | |
|-------------------|---|----------------------|
| Lab Sample ID: | JC16571-4 Date Samp | led: 03/17/16 |
| Matrix: | AIR - Ambient Air Comp. Summa ID: M111 Date Recei | ved: 03/18/16 |
| Method: | TO-15 Percent So | lids: n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | |

Report of Analysis

VOA TO15 List

| CAS No. | MW | Compound | Res | ult | RL | MDL | Units | Q | Result | RL | MDL | Units |
|-----------|--------|---------------------------|------|--------|-------|---------|-------|---|--------|------|-------|-------|
| 64-17-5 | 46.07 | Ethanol | 2.5 | | 0.50 | 0.075 | ppbv | | 4.7 | 0.94 | 0.14 | ug/m3 |
| 100-41-4 | 106.2 | Ethylbenzene | ND | | 0.20 | 0.042 | ppbv | | ND | 0.87 | 0.18 | ug/m3 |
| 141-78-6 | 88 | Ethyl Acetate | 0.2 | 1 | 0.20 | 0.075 | | | 0.76 | 0.72 | 0.27 | ug/m3 |
| 622-96-8 | 120.2 | 4-Ethyltoluene | ND | | 0.20 | 0.017 | ppbv | | ND | 0.98 | 0.084 | ug/m3 |
| 76-13-1 | 187.4 | Freon 113 | ND | | 0.20 | 0.021 | ppbv | | ND | 1.5 | 0.16 | ug/m3 |
| 76-14-2 | 170.9 | Freon 114 | ND | | 0.20 | 0.031 | | | ND | 1.4 | 0.22 | ug/m3 |
| 142-82-5 | 100.2 | Heptane | ND | | 0.20 | 0.020 | ppbv | | ND | 0.82 | 0.082 | ug/m3 |
| 87-68-3 | 260.8 | Hexachlorobutadiene | ND | | 0.20 | 0.020 | ppbv | | ND | 2.1 | 0.21 | ug/m3 |
| 110-54-3 | 86.17 | Hexane | 0.19 | 9 | 0.20 | 0.023 | ppbv | J | 0.67 | 0.70 | 0.081 | ug/m3 |
| 591-78-6 | 100 | 2-Hexanone | ND | | 0.20 | 0.045 | ppbv | | ND | 0.82 | 0.18 | ug/m3 |
| 67-63-0 | 60.1 | Isopropyl Alcohol | 0.3 | 3 | 0.20 | 0.16 | ppbv | | 0.81 | 0.49 | 0.39 | ug/m3 |
| 75-09-2 | 84.94 | Methylene chloride | 0.40 |) | 0.20 | 0.025 | ppbv | | 1.4 | 0.69 | 0.087 | ug/m3 |
| 78-93-3 | 72.11 | Methyl ethyl ketone | ND | | 0.20 | 0.048 | ppbv | | ND | 0.59 | 0.14 | ug/m3 |
| 108-10-1 | 100.2 | Methyl Isobutyl Ketone | ND | | 0.20 | 0.055 | ppbv | | ND | 0.82 | 0.23 | ug/m3 |
| 1634-04-4 | 88.15 | Methyl Tert Butyl Ether | ND | | 0.20 | 0.020 | ppbv | | ND | 0.72 | 0.072 | ug/m3 |
| 80-62-6 | 100.12 | Methylmethacrylate | ND | | 0.20 | 0.040 | ppbv | | ND | 0.82 | 0.16 | ug/m3 |
| 115-07-1 | 42 | Propylene | ND | | 0.50 | 0.032 | ppbv | | ND | 0.86 | 0.055 | ug/m3 |
| 100-42-5 | 104.1 | Styrene | ND | | 0.20 | 0.015 | ppbv | | ND | 0.85 | 0.064 | ug/m3 |
| 71-55-6 | 133.4 | 1,1,1-Trichloroethane | ND | | 0.20 | 0.024 | ppbv | | ND | 1.1 | 0.13 | ug/m3 |
| 79-34-5 | 167.9 | 1,1,2,2-Tetrachloroethane | ND | | 0.20 | 0.016 | ppbv | | ND | 1.4 | 0.11 | ug/m3 |
| 79-00-5 | 133.4 | 1,1,2-Trichloroethane | ND | | 0.20 | 0.039 | ppbv | | ND | 1.1 | 0.21 | ug/m3 |
| 120-82-1 | 181.5 | 1,2,4-Trichlorobenzene | ND | | 0.20 | 0.056 | ppbv | | ND | 1.5 | 0.42 | ug/m3 |
| 95-63-6 | 120.2 | 1,2,4-Trimethylbenzene | ND | | 0.20 | 0.015 | ppbv | | ND | 0.98 | 0.074 | ug/m3 |
| 108-67-8 | 120.2 | 1,3,5-Trimethylbenzene | ND | | 0.20 | 0.045 | ppbv | | ND | 0.98 | 0.22 | ug/m3 |
| 540-84-1 | 114.2 | 2,2,4-Trimethylpentane | ND | | 0.20 | 0.023 | ppbv | | ND | 0.93 | 0.11 | ug/m3 |
| 75-65-0 | 74.12 | Tertiary Butyl Alcohol | ND | | 0.20 | 0.053 | ppbv | | ND | 0.61 | 0.16 | ug/m3 |
| 127-18-4 | 165.8 | Tetrachloroethylene | ND | | 0.040 | 0.023 | ppbv | | ND | 0.27 | 0.16 | ug/m3 |
| 109-99-9 | 72.11 | Tetrahydrofuran | ND | | 0.20 | 0.045 | ppbv | | ND | 0.59 | 0.13 | ug/m3 |
| 108-88-3 | 92.14 | Toluene | 0.13 | 3 | 0.20 | 0.012 | ppbv | J | 0.49 | 0.75 | 0.045 | ug/m3 |
| 79-01-6 | 131.4 | Trichloroethylene | ND | | 0.040 | 0.019 | ppbv | | ND | 0.21 | 0.10 | ug/m3 |
| 75-69-4 | 137.4 | Trichlorofluoromethane | ND | | 0.20 | 0.022 | ppbv | | ND | 1.1 | 0.12 | ug/m3 |
| 75-01-4 | 62.5 | Vinyl chloride | ND | | 0.20 | 0.021 | ppbv | | ND | 0.51 | 0.054 | ug/m3 |
| 108-05-4 | 86 | Vinyl Acetate | ND | | 0.20 | 0.054 | ppbv | | ND | 0.70 | 0.19 | ug/m3 |
| | 106.2 | m,p-Xylene | ND | | 0.20 | 0.068 | ppbv | | ND | 0.87 | 0.30 | ug/m3 |
| 95-47-6 | 106.2 | o-Xylene | ND | | 0.20 | 0.051 | ppbv | | ND | 0.87 | 0.22 | ug/m3 |
| 1330-20-7 | 106.2 | Xylenes (total) | ND | | 0.20 | 0.051 | ppbv | | ND | 0.87 | 0.22 | ug/m3 |
| CAS No. | Surrog | ate Recoveries Run# | 1 | Run# : | 2 I | Limits | | | | | | |
| 460-00-4 | 4-Brom | ofluorobenzene 90% | | | e | 55-128% |) | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



JC16571

Page 2 of 2

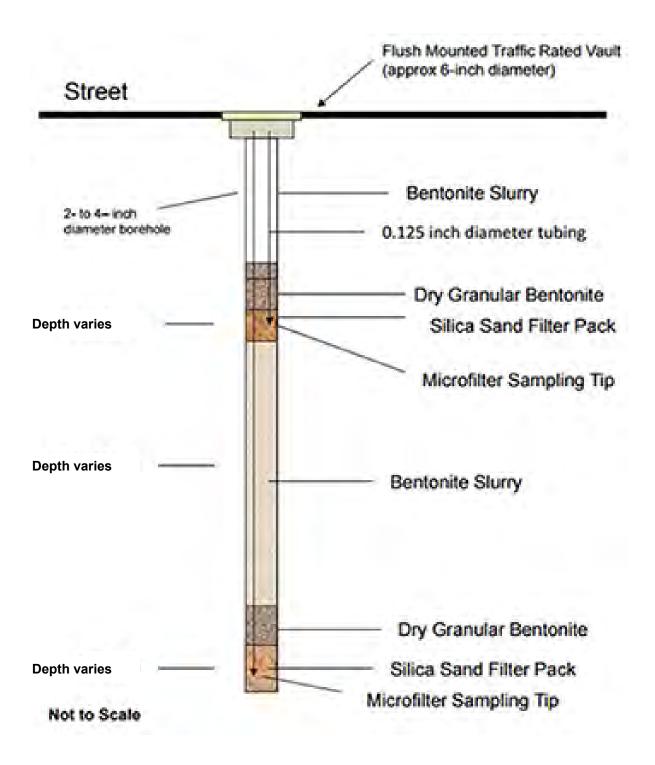
4.4 **4**

Appendix D

Typical Soil Vapor Probe and Implant Details

Remedial Investigation Report, 2002-2024 Cropsey Avenue, Brooklyn, NY March 23, 2018

Typical Soil Vapor Monitoring Probe Details



Typical Soil Vapor Probe Implants



Appendix E

Soil Data Analytical Reports



Section 4

Sample Results

Report of Analysis





| | | | - I | | J | | |
|-----------------------------|--------------|---------------|-----------------|------------|-----------|-------------------|------------------|
| Client Sa | mple ID: SB | -1-1 | | | | | |
| Lab Sam | ple ID: JC | 7512-1 | | | D | ate Sampled: 10 |)/29/15 |
| Matrix: | SO | - Soil | | | D | ate Received: 10 |)/31/15 |
| Method: | SW | /846 8260C | | | P | ercent Solids: 90 |).0 |
| Project: | 200 | 02-2024 Crops | sey Avenue, Bro | ooklyn, NY | Y | | |
| | | | - | - | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | I209229.D | 1 | 11/03/15 | SJM | n/a | n/a | VI8395 |
| Run #2 | | | | | | | |
| | Initial Weig | oht | | | | | |
| Run #1 | 5.3 g | 5 | | | | | |
| Run #2 | 5.5 6 | | | | | | |
| $\operatorname{Kull} \pi 2$ | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 36.2 | 10 | 2.3 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.52 | 0.14 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.2 | 0.32 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.1 | 0.16 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.2 | 0.25 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.2 | 0.38 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 2.0 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.1 | 0.24 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.1 | 0.24 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.1 | 0.16 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.2 | 0.50 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.1 | 0.16 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.2 | 0.27 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.1 | 0.33 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.1 | 0.57 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.1 | 0.21 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.14 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.13 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.16 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.24 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.2 | 0.38 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.15 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.14 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.62 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.82 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.62 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.1 | 0.25 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.1 | 0.12 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.1 | 0.19 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.17 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.2 | 0.47 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.2 | 1.4 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

13 of 1690

ACCUTEST JC7512

SGS

1,1,2,2-Tetrachloroethane

1,2,3-Trichlorobenzene

1,2,4-Trichlorobenzene

1,1,1-Trichloroethane

1,1,2-Trichloroethane

Trichlorofluoromethane

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Trichloroethene

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

o-Xylene

Tetrachloroethene

Toluene

79-34-5

127-18-4

108-88-3

87-61-6

120-82-1

71-55-6

79-00-5

79-01-6

75-69-4

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

Report of Analysis

| Client Samp Lab Sample Matrix: Method: Project: | | enue, Brook | lyn, NY | | Date | Sampled: Received: ent Solids: | 10/29/15 10/31/15 90.0 |
|--|---|----------------------------------|---|---|---|--------------------------------------|------------------------------|
| VOA TCL I | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 79-20-9 108-87-2 1634-04-4 108-10-1 75-09-2 100-42-5 | Isopropylbenzene Methyl Acetate Methylcyclohexane Methyl Tert Butyl Ether 4-Methyl-2-pentanone(MIBK) Methylene chloride Styrene | ND ND ND ND ND ND | 2.1 5.2 2.1 1.0 5.2 5.2 2.1 | 0.11 0.90 0.24 0.16 0.48 1.0 0.19 | ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg | | |

2.1

2.1

1.0

5.2

5.2

2.1

2.1

1.0

5.2

2.1

1.0

1.0

1.0

Run# 2

ND

18.3

ND

99%

95%

93%

93%

Run#1

0.18

0.32

0.22

0.18

0.18

0.16

0.15

0.15

0.26

0.21

0.37

0.29

0.29

Limits

70-122%

68-124%

77-125%

72-130%

ug/kg

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



14 of 1690

ACCUTEST JC7512

Page 2 of 2

4.1 4

| | | | Repo | ort of A | Analysis | | Page 1 of 3 |
|---|-------------------------------|----------------------------------|------------------------------|-----------------|---------------------------|-----------------------|------------------------------------|
| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC7 SO SW3 | 512-1 - Soil 846 8270D - S | SW846 3546 ey Avenue, Bro | ooklyn, N | Da Pe | te Received: 10 |)/29/15)/31/15).0 |
| Run #1 Run #2 | File ID 3E77583.D | DF 1 | Analyzed 11/03/15 | By AN | Prep Date 11/03/15 | Prep Batch OP88613 | Analytical Batch E3E3376 |
| Run #1 Run #2 | Initial Weig 30.8 g | ht Final V 1.0 ml | olume | | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|----------------------------|--------|-----|-----|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 72 | 27 | ug/kg | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 180 | 33 | ug/kg | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 180 | 29 | ug/kg | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 180 | 66 | ug/kg | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 180 | 160 | ug/kg | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 180 | 69 | ug/kg | |
| 95-48-7 | 2-Methylphenol | ND | 72 | 52 | ug/kg | |
| | 3&4-Methylphenol | ND | 72 | 34 | ug/kg | |
| 88-75-5 | 2-Nitrophenol | ND | 180 | 33 | ug/kg | |
| 100-02-7 | 4-Nitrophenol | ND | 360 | 61 | ug/kg | |
| 87-86-5 | Pentachlorophenol | ND | 180 | 88 | ug/kg | |
| 108-95-2 | Phenol | ND | 72 | 27 | ug/kg | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 180 | 34 | ug/kg | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 180 | 32 | ug/kg | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 180 | 29 | ug/kg | |
| 83-32-9 | Acenaphthene | ND | 36 | 34 | ug/kg | |
| 208-96-8 | Acenaphthylene | ND | 36 | 3.8 | ug/kg | |
| 98-86-2 | Acetophenone | ND | 180 | 6.1 | ug/kg | |
| 120-12-7 | Anthracene | ND | 36 | 3.1 | ug/kg | |
| 1912-24-9 | Atrazine | ND | 72 | 15 | ug/kg | |
| 56-55-3 | Benzo(a)anthracene | ND | 36 | 7.0 | ug/kg | |
| 50-32-8 | Benzo(a)pyrene | ND | 36 | 7.7 | ug/kg | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 36 | 7.4 | ug/kg | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 36 | 11 | ug/kg | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 36 | 8.0 | ug/kg | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 72 | 8.2 | ug/kg | |
| 85-68-7 | Butyl benzyl phthalate | ND | 72 | 19 | ug/kg | |
| 92-52-4 | 1,1'-Biphenyl | ND | 72 | 6.7 | ug/kg | |
| 100-52-7 | Benzaldehyde | ND | 180 | 9.0 | ug/kg | |
| 91-58-7 | 2-Chloronaphthalene | ND | 72 | 5.2 | ug/kg | |
| 106-47-8 | 4-Chloroaniline | ND | 180 | 9.6 | ug/kg | |
| 86-74-8 | Carbazole | ND | 72 | 4.0 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





15 of 1690

ACCUTEST JC7512

Report of Analysis

| Client Sample ID: | SB-1-1 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC7512-1 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 90.0 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|--------|------|-------|---|
| 105-60-2 | Caprolactam | ND | 72 | 23 | ug/kg | |
| 218-01-9 | Chrysene | ND | 36 | 5.8 | ug/kg | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 72 | 8.2 | ug/kg | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 72 | 15 | ug/kg | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 72 | 8.3 | ug/kg | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 72 | 6.8 | ug/kg | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 36 | 6.8 | ug/kg | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 36 | 9.3 | ug/kg | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 72 | 24 | ug/kg | |
| 123-91-1 | 1,4-Dioxane | ND | 36 | 24 | ug/kg | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 36 | 13 | ug/kg | |
| 132-64-9 | Dibenzofuran | ND | 72 | 5.0 | ug/kg | |
| 84-74-2 | Di-n-butyl phthalate | ND | 72 | 4.3 | ug/kg | |
| 117-84-0 | Di-n-octyl phthalate | ND | 72 | 4.9 | ug/kg | |
| 84-66-2 | Diethyl phthalate | ND | 72 | 4.6 | ug/kg | |
| 131-11-3 | Dimethyl phthalate | ND | 72 | 5.2 | ug/kg | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 223 | 72 | 13 | ug/kg | |
| 206-44-0 | Fluoranthene | ND | 36 | 4.4 | ug/kg | |
| 86-73-7 | Fluorene | ND | 36 | 4.3 | ug/kg | |
| 118-74-1 | Hexachlorobenzene | ND | 72 | 7.1 | ug/kg | |
| 87-68-3 | Hexachlorobutadiene | ND | 36 | 9.6 | ug/kg | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 360 | 57 | ug/kg | |
| 67-72-1 | Hexachloroethane | ND | 180 | 12 | ug/kg | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 36 | 19 | ug/kg | |
| 78-59-1 | Isophorone | ND | 72 | 6.7 | ug/kg | |
| 91-57-6 | 2-Methylnaphthalene | ND | 72 | 6.7 | ug/kg | |
| 88-74-4 | 2-Nitroaniline | ND | 180 | 8.2 | ug/kg | |
| 99-09-2 | 3-Nitroaniline | ND | 180 | 10 | ug/kg | |
| 100-01-6 | 4-Nitroaniline | ND | 180 | 12 | ug/kg | |
| 91-20-3 | Naphthalene | ND | 36 | 5.8 | ug/kg | |
| 98-95-3 | Nitrobenzene | ND | 72 | 11 | ug/kg | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 72 | 11 | ug/kg | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 180 | 19 | ug/kg | |
| 85-01-8 | Phenanthrene | ND | 36 | 4.0 | ug/kg | |
| 129-00-0 | Pyrene | ND | 36 | 4.5 | ug/kg | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 180 | 8.7 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 367-12-4 | 2-Fluorophenol | 80% | | 30-1 | 06% | |
| | | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

4.1

Page 2 of 3



16 of 1690

ACCUTEST JC7512

Report of Analysis

| Client Sample ID: | SB-1-1 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-1 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 90.0 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 4165-62-2 | Phenol-d5 | 77% | | 30-106% |
| 118-79-6 | 2,4,6-Tribromophenol | 93% | | 24-140% |
| 4165-60-0 | Nitrobenzene-d5 | 106% | | 26-122% |
| 321-60-8 | 2-Fluorobiphenyl | 90% | | 36-112% |
| 1718-51-0 | Terphenyl-d14 | 90% | | 36-132% |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



17 of 1690

ACCUTEST JC7512

4.1 **4**

Page 3 of 3

| | | | Repo | ort of A | nalysis | | Page 1 of |
|--|-----------------------------------|--------------------------|------------------------------|------------------|---------------------------|-----------------------|-----------------------------|
| Client San Lab Samp Matrix: Method: Project: | le ID: JC7512 SO - So SW846 | oil 8151 SW | 7846 3550C ey Avenue, Bro | ooklyn, NY | Da Pe | ate Received: 10 |)/29/15)/31/15).0 |
| Run #1 Run #2 | File ID OA112008.D | DF 1 | Analyzed 11/05/15 | By VDT | Prep Date 11/03/15 | Prep Batch OP88618 | Analytical Batch GOA3882 |
| Run #1 Run #2 | Initial Weight 15.2 g | Final V 5.0 ml | olume | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|------------------|-------------------|-------------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 18 3.7 3.7 | 8.1 1.1 1.4 | ug/kg ug/kg ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2,4-DCAA | 23% 20% | | 10-1 10-1 | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



18 of 1690 ACCUTEST JC7512

4

| Report of Analysis PA | | | | | | | | |
|---|------------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|-----------------------|------------------------------|--|
| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC7512 SO - So SW846 | 2-1 oil 5 8081B | SW846 3546 sey Avenue, Bro | ooklyn, N | I F | Date Received: | 10/29/15 10/31/15 90.0 | |
| Run #1 Run #2 | File ID 1G116787.D | DF 1 | Analyzed 11/05/15 | By RK | Prep Date 11/03/15 | Prep Batch OP88620 | Analytical Batch G1G3831 | |
| Run #1 Run #2 | Initial Weight 15.6 g | Final V 10.0 m | | | | | | |

Pesticide TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|------------------------|--------|--------|------|-------|---|
| 309-00-2 | Aldrin | ND | 0.71 | 0.64 | ug/kg | |
| 319-84-6 | alpha-BHC | ND | 0.71 | 0.48 | ug/kg | |
| 319-85-7 | beta-BHC | ND | 0.71 | 0.44 | ug/kg | |
| 319-86-8 | delta-BHC | ND | 0.71 | 0.28 | ug/kg | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.71 | 0.32 | ug/kg | |
| 5103-71-9 | alpha-Chlordane | ND | 0.71 | 0.38 | ug/kg | |
| 5103-74-2 | gamma-Chlordane | ND | 0.71 | 0.54 | ug/kg | |
| 60-57-1 | Dieldrin | ND | 0.71 | 0.56 | ug/kg | |
| 72-54-8 | 4,4' -DDD ^a | 0.72 | 0.71 | 0.26 | ug/kg | |
| 72-55-9 | 4,4'-DDE | ND | 0.71 | 0.24 | ug/kg | |
| 50-29-3 | 4,4' -DDT | 1.9 | 0.71 | 0.27 | ug/kg | |
| 72-20-8 | Endrin | ND | 0.71 | 0.25 | ug/kg | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.71 | 0.41 | ug/kg | |
| 7421-93-4 | Endrin aldehyde | ND | 0.71 | 0.53 | ug/kg | |
| 959-98-8 | Endosulfan-I | ND | 0.71 | 0.23 | ug/kg | |
| 33213-65-9 | Endosulfan-II | ND | 0.71 | 0.67 | ug/kg | |
| 76-44-8 | Heptachlor | ND | 0.71 | 0.59 | ug/kg | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.71 | 0.29 | ug/kg | |
| 72-43-5 | Methoxychlor | ND | 1.4 | 0.40 | ug/kg | |
| 53494-70-5 | Endrin ketone | ND | 0.71 | 0.37 | ug/kg | |
| 8001-35-2 | Toxaphene | ND | 18 | 12 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 877-09-8 | Tetrachloro-m-xylene | 80% | | 24-1 | 36% | |
| 877-09-8 | Tetrachloro-m-xylene | 78% | | 24-1 | 36% | |
| 2051-24-3 | Decachlorobiphenyl | 67% | | 10-1 | 53% | |
| 2051-24-3 | Decachlorobiphenyl | 95% | | 10-1 | 53% | |

(a) More than 40 % RPD for detected concentrations between the two GC columns.

- ND = Not detected MDL = Method Detection Limit
- RL = Reporting Limit
- E = Indicates value exceeds calibration range
- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4.1

JC7512

| | Report of Analysis Page 1 of 1 | | | | | | |
|--|---------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|-----------------------|------------------------------|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC751 SO - S SW84 | 2-1 oil 6 8082A | SW846 3546 sey Avenue, Bro | ooklyn, N | D P | ate Received: | 10/29/15 10/31/15 90.0 |
| Run #1 Run #2 | File ID XX179842.D | DF 1 | Analyzed 11/06/15 | By JR | Prep Date 11/03/15 | Prep Batch OP88619 | Analytical Batch GXX5510 |
| Run #1 Run #2 | Initial Weight 15.0 g | Final V 10.0 m | 0-0 | | | | |

PCB List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|--|--|-------------------------------------|--|----------------------------------|---|---|
| 12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 | Aroclor 1016 Aroclor 1221 Aroclor 1232 Aroclor 1242 Aroclor 1248 Aroclor 1254 | ND ND ND ND ND | 37 37 37 37 37 37 37 | 12 22 12 17 11 16 | ug/kg ug/kg ug/kg ug/kg ug/kg | |
| 11096-82-5 11100-14-4 37324-23-5 | Aroclor 1260 Aroclor 1268 Aroclor 1262 | ND ND ND | 37 37 37 | 16 11 10 | ug/kg ug/kg ug/kg | |
| CAS No. 877-09-8 877-09-8 2051-24-3 2051-24-3 | Surrogate Recoveries Tetrachloro-m-xylene Tetrachloro-m-xylene Decachlorobiphenyl Decachlorobiphenyl | Run# 1 100% 99% 94% 96% | Run# 2 | 20-1 20-1 12-1 12-1 | 52% 52% 57% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| Client Sample ID: | SB-1-1 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC7512-1 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| | | Percent Solids: | 90.0 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|---------|-------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | 6060 | 55 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Antimony | < 2.2 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Arsenic | 2.2 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Barium | 25.2 | 22 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Beryllium | 0.44 | 0.22 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Cadmium | < 0.55 | 0.55 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Calcium | 1080 | 550 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Chromium | 18.6 | 1.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Cobalt | 9.4 | 5.5 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Copper | 10.2 | 2.8 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Iron | 12700 | 55 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Lead | 5.6 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Magnesium | 2800 | 550 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Manganese | 272 | 1.7 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Mercury | < 0.035 | 0.035 | mg/kg | 1 | 11/04/15 | 11/04/15 ма | SW846 7471B ¹ | SW846 7471B ⁴ |
| Nickel | 48.6 | 4.4 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Potassium | 1300 | 1100 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Selenium | < 2.2 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Silver | 0.64 | 0.55 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Sodium | < 1100 | 1100 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Thallium | < 1.1 | 1.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Vanadium | 20.7 | 5.5 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Zinc | 24.0 | 5.5 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |

(1) Instrument QC Batch: MA37945

(2) Instrument QC Batch: MA37978

(3) Prep QC Batch: MP90046

(4) Prep QC Batch: MP90074



Page 1 of 1



| | | | - I | | | | |
|-----------|-------------|---------------|-----------------|------------|-----------|-------------------|------------------|
| Client Sa | mple ID: SB | -1-5 | | | | | |
| Lab Sam | ple ID: JC | 7512-2 | | | D | ate Sampled: 10 |)/29/15 |
| Matrix: | SO | - Soil | | | D | ate Received: 10 |)/31/15 |
| Method: | SW | /846 8260C | | | Р | ercent Solids: 96 | 5.1 |
| Project: | 200 | 02-2024 Crops | sey Avenue, Bro | ooklyn, NY | Y | | |
| | | | - | - | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | I209230.D | 1 | 11/03/15 | SJM | n/a | n/a | VI8395 |
| Run #2 | | | | | | | |
| | Initial Wei | oht | | | | | |
| Run #1 | 4.6 g | 5 | | | | | |
| Run #2 | 1.0 5 | | | | | | |
| itun 112 | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 11 | 2.5 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.57 | 0.15 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.7 | 0.35 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.3 | 0.18 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.7 | 0.27 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.7 | 0.41 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 11 | 2.2 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.3 | 0.26 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.3 | 0.26 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.3 | 0.18 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.7 | 0.54 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.3 | 0.17 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.7 | 0.30 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.3 | 0.36 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.3 | 0.62 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.3 | 0.23 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.1 | 0.15 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.1 | 0.14 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.1 | 0.18 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.1 | 0.25 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.7 | 0.41 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.1 | 0.16 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.1 | 0.15 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.1 | 0.67 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.1 | 0.88 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.1 | 0.67 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.3 | 0.27 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.3 | 0.13 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.3 | 0.20 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.1 | 0.18 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.7 | 0.51 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.7 | 1.5 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

 $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



Page 1 of 2

SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Report of | f Analysis |
|------------------|------------|
|------------------|------------|

| Client Samp Lab Sample Matrix: Method: Project: | | enue, Brool | clyn, NY | | Date | Sampled: Received: ent Solids: | 10/29/15 10/31/15 96.1 |
|---|----------------------------|-------------|----------|------|-------|--------------------------------------|------------------------------|
| VOA TCL I | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 2.3 | 0.12 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 5.7 | 0.97 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 2.3 | 0.26 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.1 | 0.17 | ug/kg | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.7 | 0.52 | ug/kg | | |
| 75-09-2 | Methylene chloride | ND | 5.7 | 1.1 | ug/kg | | |
| 100-42-5 | Styrene | ND | 2.3 | 0.20 | ug/kg | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.3 | 0.20 | ug/kg | | |
| 127-18-4 | Tetrachloroethene | 19.6 | 2.3 | 0.34 | ug/kg | | |
| 108-88-3 | Toluene | ND | 1.1 | 0.24 | ug/kg | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.7 | 0.20 | ug/kg | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.7 | 0.19 | ug/kg | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.3 | 0.17 | ug/kg | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.3 | 0.17 | ug/kg | | |
| 79-01-6 | Trichloroethene | 0.40 | 1.1 | 0.17 | ug/kg | J | |
| 75-69-4 | Trichlorofluoromethane | ND | 5.7 | 0.28 | ug/kg | | |
| 75-01-4 | Vinyl chloride | ND | 2.3 | 0.22 | ug/kg | | |
| | m,p-Xylene | ND | 1.1 | 0.40 | ug/kg | | |
| 95-47-6 | o-Xylene | ND | 1.1 | 0.31 | ug/kg | | |
| 1330-20-7 | Xylene (total) | ND | 1.1 | 0.31 | ug/kg | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 1868-53-7 | Dibromofluoromethane | 94% | | 70-1 | 22% | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 93% | | 68-1 | | | |
| 2037-26-5 | Toluene-D8 | 93% | | 77-1 | | | |
| 460-00-4 | 4-Bromofluorobenzene | 89% | | 72-1 | | | |

| ND = Not detected | MDL = Method Detection Limit | | | |
|---|------------------------------|--|--|--|
| RL = Reporting Limit | | | | |
| E = Indicates value exceeds calibration range | | | | |

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



Page 2 of 2



| | | | Repo | ort of A | Analysis | | Page 1 of 3 |
|---|---------------------------------|-----------------------|-------------------------------|-----------------|---------------------------|-----------------------|------------------------------------|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC751 SO - S SW84 | 2-2 oil 6 8270D | SW846 3546 sey Avenue, Bro | ooklyn, N | Da Pe | L . |)/29/15)/31/15 5.1 |
| Run #1 Run #2 | File ID 3E77584.D | DF 1 | Analyzed 11/03/15 | By AN | Prep Date 11/03/15 | Prep Batch OP88613 | Analytical Batch E3E3376 |
| Run #1 Run #2 | Initial Weight 30.6 g | Final V 1.0 ml | olume | | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units Q | |
|-----------|----------------------------|--------|-----|-----|---------|--|
| 95-57-8 | 2-Chlorophenol | ND | 68 | 25 | ug/kg | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 170 | 31 | ug/kg | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 170 | 27 | ug/kg | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 170 | 62 | ug/kg | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 170 | 150 | ug/kg | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 170 | 65 | ug/kg | |
| 95-48-7 | 2-Methylphenol | ND | 68 | 49 | ug/kg | |
| | 3&4-Methylphenol | ND | 68 | 33 | ug/kg | |
| 88-75-5 | 2-Nitrophenol | ND | 170 | 31 | ug/kg | |
| 100-02-7 | 4-Nitrophenol | ND | 340 | 58 | ug/kg | |
| 87-86-5 | Pentachlorophenol | ND | 170 | 83 | ug/kg | |
| 108-95-2 | Phenol | ND | 68 | 25 | ug/kg | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 170 | 32 | ug/kg | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 170 | 31 | ug/kg | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 170 | 27 | ug/kg | |
| 83-32-9 | Acenaphthene | ND | 34 | 32 | ug/kg | |
| 208-96-8 | Acenaphthylene | ND | 34 | 3.6 | ug/kg | |
| 98-86-2 | Acetophenone | ND | 170 | 5.8 | ug/kg | |
| 120-12-7 | Anthracene | ND | 34 | 2.9 | ug/kg | |
| 1912-24-9 | Atrazine | ND | 68 | 14 | ug/kg | |
| 56-55-3 | Benzo(a)anthracene | ND | 34 | 6.6 | ug/kg | |
| 50-32-8 | Benzo(a)pyrene | ND | 34 | 7.2 | ug/kg | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 34 | 7.0 | ug/kg | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 34 | 10 | ug/kg | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 34 | 7.6 | ug/kg | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 68 | 7.8 | ug/kg | |
| 85-68-7 | Butyl benzyl phthalate | ND | 68 | 18 | ug/kg | |
| 92-52-4 | 1,1'-Biphenyl | ND | 68 | 6.3 | ug/kg | |
| 100-52-7 | Benzaldehyde | ND | 170 | 8.5 | ug/kg | |
| 91-58-7 | 2-Chloronaphthalene | ND | 68 | 4.9 | ug/kg | |
| 106-47-8 | 4-Chloroaniline | ND | 170 | 9.0 | ug/kg | |
| 86-74-8 | Carbazole | ND | 68 | 3.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

```
J = Indicates an estimated value
```

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





| Client Sample ID: | SB-1-5 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-2 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 96.1 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|--------|-------|-------|---|
| 105-60-2 | Caprolactam | ND | 68 | 22 | ug/kg | |
| 218-01-9 | Chrysene | ND | 34 | 5.5 | ug/kg | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 68 | 7.7 | ug/kg | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 68 | 14 | ug/kg | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 68 | 7.8 | ug/kg | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 68 | 6.4 | ug/kg | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 34 | 6.4 | ug/kg | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 34 | 8.8 | ug/kg | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 68 | 22 | ug/kg | |
| 123-91-1 | 1,4-Dioxane | ND | 34 | 23 | ug/kg | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 34 | 12 | ug/kg | |
| 132-64-9 | Dibenzofuran | ND | 68 | 4.7 | ug/kg | |
| 84-74-2 | Di-n-butyl phthalate | ND | 68 | 4.0 | ug/kg | |
| 117-84-0 | Di-n-octyl phthalate | ND | 68 | 4.6 | ug/kg | |
| 84-66-2 | Diethyl phthalate | ND | 68 | 4.3 | ug/kg | |
| 131-11-3 | Dimethyl phthalate | ND | 68 | 4.9 | ug/kg | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 68 | 12 | ug/kg | |
| 206-44-0 | Fluoranthene | ND | 34 | 4.1 | ug/kg | |
| 86-73-7 | Fluorene | ND | 34 | 4.0 | ug/kg | |
| 118-74-1 | Hexachlorobenzene | ND | 68 | 6.7 | ug/kg | |
| 87-68-3 | Hexachlorobutadiene | ND | 34 | 9.0 | ug/kg | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 340 | 54 | ug/kg | |
| 67-72-1 | Hexachloroethane | ND | 170 | 11 | ug/kg | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 34 | 18 | ug/kg | |
| 78-59-1 | Isophorone | ND | 68 | 6.4 | ug/kg | |
| 91-57-6 | 2-Methylnaphthalene | ND | 68 | 6.4 | ug/kg | |
| 88-74-4 | 2-Nitroaniline | ND | 170 | 7.7 | ug/kg | |
| 99-09-2 | 3-Nitroaniline | ND | 170 | 9.7 | ug/kg | |
| 100-01-6 | 4-Nitroaniline | ND | 170 | 11 | ug/kg | |
| 91-20-3 | Naphthalene | ND | 34 | 5.4 | ug/kg | |
| 98-95-3 | Nitrobenzene | ND | 68 | 11 | ug/kg | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 68 | 10 | ug/kg | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 170 | 18 | ug/kg | |
| 85-01-8 | Phenanthrene | ND | 34 | 3.8 | ug/kg | |
| 129-00-0 | Pyrene | ND | 34 | 4.3 | ug/kg | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 170 | 8.2 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 367-12-4 | 2-Fluorophenol | 78% | | 30-10 | 06% | |
| | | | | | | |

ND = Not detectedMDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 3

25 of 1690

ACCUTEST JC7512

Report of Analysis

| Client Sample ID: | SB-1-5 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC7512-2 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 96.1 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | · · · | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 4165-62-2 | Phenol-d5 | 76% | | 30-106% |
| 118-79-6 | 2,4,6-Tribromophenol | 91% | | 24-140% |
| 4165-60-0 | Nitrobenzene-d5 | 102% | | 26-122% |
| 321-60-8 | 2-Fluorobiphenyl | 90% | | 36-112% |
| 1718-51-0 | Terphenyl-d14 | 86% | | 36-132% |

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



26 of 1690

ACCUTEST JC7512

Page 3 of 3



| | | | Repo | ort of A | nalysis | | Page 1 of |
|--|------------------------------------|-------------------|------------------------------|------------------|---------------------------|-----------------------|------------------------------------|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: JC7512 SO - So SW846 | oil 8151 SV | W846 3550C ey Avenue, Bro | ooklyn, N | Da Pe | ate Received: 10 | 0/29/15 0/31/15 5.1 |
| Run #1 Run #2 | File ID OA112009.D | DF 2 | Analyzed 11/05/15 | By VDT | Prep Date 11/03/15 | Prep Batch OP88618 | Analytical Batch GOA3882 |
| Run #1 Run #2 | Initial Weight 15.2 g | Final V 5.0 ml | olume | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|------------------|------------------|-------------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 34 6.8 6.8 | 15 2.0 2.5 | ug/kg ug/kg ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2,4-DCAA | 40% 31% | | 10-1 10-1 | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



27 of 1690 ACCUTEST JC7512



| | | | Repo | ort of A | Analysis | | Page 1 of 1 |
|--|----------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|-----------------------|------------------------------------|
| Client San Lab Samp Matrix: Method: Project: | ole ID: JC751 SO - S SW840 | 2-2 oil 5 8081B | SW846 3546 sey Avenue, Bro | ooklyn, N | Da Pe | L |)/29/15)/31/15 5.1 |
| Run #1 Run #2 | File ID 1G116788.D | DF 1 | Analyzed 11/05/15 | By RK | Prep Date 11/03/15 | Prep Batch OP88620 | Analytical Batch G1G3831 |
| Run #1 Run #2 | Initial Weight 15.7 g | Final V 10.0 m | | | | | |

Pesticide TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------|--------|--------|------|-------|---|
| 309-00-2 | Aldrin | ND | 0.66 | 0.59 | ug/kg | |
| 319-84-6 | alpha-BHC | ND | 0.66 | 0.44 | ug/kg | |
| 319-85-7 | beta-BHC | ND | 0.66 | 0.41 | ug/kg | |
| 319-86-8 | delta-BHC | ND | 0.66 | 0.26 | ug/kg | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.66 | 0.30 | ug/kg | |
| 5103-71-9 | alpha-Chlordane | ND | 0.66 | 0.35 | ug/kg | |
| 5103-74-2 | gamma-Chlordane | ND | 0.66 | 0.51 | ug/kg | |
| 60-57-1 | Dieldrin | ND | 0.66 | 0.52 | ug/kg | |
| 72-54-8 | 4,4'-DDD | ND | 0.66 | 0.25 | ug/kg | |
| 72-55-9 | 4,4'-DDE | ND | 0.66 | 0.22 | ug/kg | |
| 50-29-3 | 4,4'-DDT | ND | 0.66 | 0.25 | ug/kg | |
| 72-20-8 | Endrin | ND | 0.66 | 0.23 | ug/kg | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.66 | 0.38 | ug/kg | |
| 7421-93-4 | Endrin aldehyde | ND | 0.66 | 0.49 | ug/kg | |
| 959-98-8 | Endosulfan-I | ND | 0.66 | 0.22 | ug/kg | |
| 33213-65-9 | Endosulfan-II | ND | 0.66 | 0.63 | ug/kg | |
| 76-44-8 | Heptachlor | ND | 0.66 | 0.54 | ug/kg | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.66 | 0.27 | ug/kg | |
| 72-43-5 | Methoxychlor | ND | 1.3 | 0.37 | ug/kg | |
| 53494-70-5 | Endrin ketone | ND | 0.66 | 0.35 | ug/kg | |
| 8001-35-2 | Toxaphene | ND | 17 | 11 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 877-09-8 | Tetrachloro-m-xylene | 68% | | 24-1 | 36% | |
| 877-09-8 | Tetrachloro-m-xylene | 73% | | 24-1 | 36% | |
| 2051-24-3 | Decachlorobiphenyl | 50% | | 10-1 | 53% | |
| 2051-24-3 | Decachlorobiphenyl | 86% | | 10-1 | 53% | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

- E = Indicates value exceeds calibration range
- J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





| | | | Repo | ort of A | Analysis | | Page 1 of 1 |
|--|------------------------------------|-----------------------|-------------------------------|-----------|-----------|------------|---------------------------|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC7512 SO - So SW846 | 2-2 pil 5 8082A | SW846 3546 sey Avenue, Bro | ooklyn, N | D P | L | 0/29/15 0/31/15 6.1 |
| D #1 | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 Run #2 | XX179843.D | 1 | 11/06/15 | JR | 11/03/15 | OP88619 | GXX5510 |
| D /// | Initial Weight | | olume | | | | |
| Run #1 Run #2 | 15.7 g | 10.0 m | 1 | | | | |

PCB List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|--|--|--|--|--|---|---|
| 12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 11096-82-5 11100-14-4 | Aroclor 1016 Aroclor 1221 Aroclor 1232 Aroclor 1242 Aroclor 1248 Aroclor 1254 Aroclor 1260 Aroclor 1268 | ND ND ND ND ND ND ND | 33 33 33 33 33 33 33 33 33 33 | 11 19 11 15 10 15 14 10 | ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg | |
| 37324-23-5 CAS No. | Aroclor 1262 Surrogate Recoveries | ND Run# 1 | 33 Run# 2 | 9.3 Limi | ug/kg its | |
| 877-09-8 877-09-8 2051-24-3 2051-24-3 | Tetrachloro-m-xylene Tetrachloro-m-xylene Decachlorobiphenyl Decachlorobiphenyl | 93% 93% 84% 86% | | 20-1 20-1 12-1 12-1 | 52% 57% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| Client Sample ID: | SB-1-5 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-2 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| | | Percent Solids: | 96.1 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|---------|-------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | 4560 | 50 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Antimony | < 2.0 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Arsenic | < 2.0 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Barium | 23.7 | 20 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Beryllium | 0.29 | 0.20 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Cadmium | < 0.50 | 0.50 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Calcium | 1020 | 500 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Chromium | 12.5 | 0.99 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Cobalt | 5.3 | 5.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Copper | 9.3 | 2.5 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Iron | 9740 | 50 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Lead | 4.2 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Magnesium | 2450 | 500 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Manganese | 244 | 1.5 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Mercury | < 0.032 | 0.032 | mg/kg | 1 | 11/04/15 | 11/04/15 ма | SW846 7471B ¹ | SW846 7471B ⁴ |
| Nickel | 50.7 | 4.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Potassium | < 990 | 990 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Selenium | < 2.0 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Silver | < 0.50 | 0.50 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Sodium | < 990 | 990 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Thallium | < 0.99 | 0.99 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Vanadium | 16.1 | 5.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Zinc | 20.8 | 5.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |

(1) Instrument QC Batch: MA37945

(2) Instrument QC Batch: MA37978

(3) Prep QC Batch: MP90046

(4) Prep QC Batch: MP90074



Page 1 of 1



| | | | | | | | ruge r or 2 |
|-----------------|--------------|--------------|----------------|-----------|-----------|----------------|------------------|
| Client Sar | nple ID: SB- | -2-1 | | | | | |
| Lab Samp | le ID: JC7 | 7512-3 | | | D | ate Sampled: | 10/29/15 |
| Matrix: | SO | - Soil | | | D | Date Received: | 10/31/15 |
| Method: | SW | '846 8260C | | | Р | ercent Solids: | 87.6 |
| Project: | 200 | 2-2024 Crops | ey Avenue, Bro | oklyn, NY | | | |
| | | | - | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | I209359.D | 1 | 11/06/15 | SJM | n/a | n/a | VI8400 |
| Run #2 | | | | | | | |
| | | | | | | | |
| | Initial Weig | ght | | | | | |
| Run #1 | 5.4 g | | | | | | |
| Run #2 | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 75.1 | 11 | 2.4 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.53 | 0.14 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.3 | 0.33 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.1 | 0.16 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.3 | 0.25 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.3 | 0.38 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 11 | 2.0 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.1 | 0.24 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.1 | 0.24 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.1 | 0.16 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.3 | 0.51 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.1 | 0.16 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.3 | 0.28 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.1 | 0.33 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.1 | 0.58 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.1 | 0.22 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.1 | 0.14 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.1 | 0.13 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.1 | 0.17 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.1 | 0.24 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.3 | 0.38 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.1 | 0.15 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.1 | 0.14 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.1 | 0.63 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.1 | 0.82 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.1 | 0.63 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.1 | 0.25 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.1 | 0.12 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.1 | 0.19 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.1 | 0.17 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.3 | 0.47 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.3 | 1.4 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

 $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



31 of 1690

ACCUTEST JC7512

Page 1 of 2

E = Indicates value exceeds calibration range

J = Indicates an estimated value

1,1,2,2-Tetrachloroethane

1,2,3-Trichlorobenzene

1,2,4-Trichlorobenzene

1,1,1-Trichloroethane

1,1,2-Trichloroethane

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Trichloroethene

m,p-Xylene

Xylene (total)

Toluene-D8

o-Xylene

Tetrachloroethene

Toluene

Report of Analysis

| Client Sam | ple ID: SB-2-1 | | | | | | |
|------------|----------------------------|------------|----------|------|-------|------------------|----------|
| Lab Sample | ID: JC7512-3 | | | | Date | Sampled: | 10/29/15 |
| Matrix: | SO - Soil | | | | Date | Received: | 10/31/15 |
| Method: | SW846 8260C | | | | Perce | ent Solids: | 87.6 |
| Project: | 2002-2024 Cropsey Av | enue, Broo | klyn, NY | | | | |
| | | | | | | | |
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | 0 | |
| CAS NO. | Compound | Result | KL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 2.1 | 0.11 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 5.3 | 0.91 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 2.1 | 0.24 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.1 | 0.16 | ug/kg | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.3 | 0.49 | ug/kg | | |
| 75-09-2 | Methylene chloride | ND | 5.3 | 1.0 | ug/kg | | |
| 100-42-5 | Styrene | ND | 2.1 | 0.19 | ug/kg | | |

2.1

2.1

1.1

5.3

5.3

2.1

2.1

1.1

5.3

2.1

1.1

1.1

1.1

Run# 2

0.18

0.32

0.22

0.19

0.18

0.16

0.16

0.16

0.26

0.21

0.37

0.29

0.29

Limits

70-122%

68-124%

77-125%

72-130%

ug/kg

J

ND

34.7

ND

ND

ND

ND

ND

0.70

ND

ND

ND

ND

ND

94%

95%

94%

98%

Run#1

75-69-4 Trichlorofluoromethane 75-01-4 Vinyl chloride

79-34-5

127-18-4

108-88-3

87-61-6

120-82-1

71-55-6

79-00-5

79-01-6

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

| ND = Not detected | MDL = Method Detection Limit |
|-------------------------|------------------------------|
| RL = Reporting Limit | |
| E = Indicates value exc | eeds calibration range |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



| | Report of Analysis | | | | | | | |
|---|---------------------------------|-----------------------|------------------------------|-----------------|---------------------------|-----------------------|------------------------------------|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC751 SO - S SW84 | 2-3 oil 6 8270D | SW846 3546 ey Avenue, Bro | ooklyn, N | Da Pe | te Received: 10 |)/29/15)/31/15 7.6 | |
| Run #1 Run #2 | File ID 3E77585.D | DF 1 | Analyzed 11/03/15 | By AN | Prep Date 11/03/15 | Prep Batch OP88613 | Analytical Batch E3E3376 | |
| Run #1 Run #2 | Initial Weight 31.6 g | Final V 1.0 ml | olume | | | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units Q | |
|-----------|----------------------------|--------|-----|-----|---------|--|
| 95-57-8 | 2-Chlorophenol | ND | 72 | 27 | ug/kg | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 180 | 33 | ug/kg | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 180 | 29 | ug/kg | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 180 | 66 | ug/kg | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 180 | 160 | ug/kg | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 180 | 69 | ug/kg | |
| 95-48-7 | 2-Methylphenol | ND | 72 | 52 | ug/kg | |
| | 3&4-Methylphenol | ND | 72 | 35 | ug/kg | |
| 88-75-5 | 2-Nitrophenol | ND | 180 | 33 | ug/kg | |
| 100-02-7 | 4-Nitrophenol | ND | 360 | 61 | ug/kg | |
| 87-86-5 | Pentachlorophenol | ND | 180 | 88 | ug/kg | |
| 108-95-2 | Phenol | ND | 72 | 27 | ug/kg | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 180 | 34 | ug/kg | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 180 | 33 | ug/kg | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 180 | 29 | ug/kg | |
| 83-32-9 | Acenaphthene | ND | 36 | 34 | ug/kg | |
| 208-96-8 | Acenaphthylene | ND | 36 | 3.8 | ug/kg | |
| 98-86-2 | Acetophenone | ND | 180 | 6.1 | ug/kg | |
| 120-12-7 | Anthracene | ND | 36 | 3.1 | ug/kg | |
| 1912-24-9 | Atrazine | ND | 72 | 15 | ug/kg | |
| 56-55-3 | Benzo(a)anthracene | ND | 36 | 7.0 | ug/kg | |
| 50-32-8 | Benzo(a)pyrene | ND | 36 | 7.7 | ug/kg | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 36 | 7.4 | ug/kg | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 36 | 11 | ug/kg | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 36 | 8.1 | ug/kg | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 72 | 8.2 | ug/kg | |
| 85-68-7 | Butyl benzyl phthalate | ND | 72 | 20 | ug/kg | |
| 92-52-4 | 1,1'-Biphenyl | ND | 72 | 6.7 | ug/kg | |
| 100-52-7 | Benzaldehyde | ND | 180 | 9.0 | ug/kg | |
| 91-58-7 | 2-Chloronaphthalene | ND | 72 | 5.2 | ug/kg | |
| 106-47-8 | 4-Chloroaniline | ND | 180 | 9.6 | ug/kg | |
| 86-74-8 | Carbazole | ND | 72 | 4.0 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

```
J = Indicates an estimated value
```

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





| Client Sample ID: | SB-2-1 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-3 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 87.6 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|--------|------|-------|---|
| 105-60-2 | Caprolactam | ND | 72 | 23 | ug/kg | |
| 218-01-9 | Chrysene | ND | 36 | 5.8 | ug/kg | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 72 | 8.2 | ug/kg | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 72 | 15 | ug/kg | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 72 | 8.3 | ug/kg | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 72 | 6.8 | ug/kg | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 36 | 6.8 | ug/kg | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 36 | 9.3 | ug/kg | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 72 | 24 | ug/kg | |
| 123-91-1 | 1,4-Dioxane | ND | 36 | 24 | ug/kg | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 36 | 13 | ug/kg | |
| 132-64-9 | Dibenzofuran | ND | 72 | 5.0 | ug/kg | |
| 84-74-2 | Di-n-butyl phthalate | ND | 72 | 4.3 | ug/kg | |
| 117-84-0 | Di-n-octyl phthalate | ND | 72 | 4.9 | ug/kg | |
| 84-66-2 | Diethyl phthalate | ND | 72 | 4.6 | ug/kg | |
| 131-11-3 | Dimethyl phthalate | ND | 72 | 5.2 | ug/kg | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 1680 | 72 | 13 | ug/kg | |
| 206-44-0 | Fluoranthene | ND | 36 | 4.4 | ug/kg | |
| 86-73-7 | Fluorene | ND | 36 | 4.3 | ug/kg | |
| 118-74-1 | Hexachlorobenzene | ND | 72 | 7.1 | ug/kg | |
| 87-68-3 | Hexachlorobutadiene | ND | 36 | 9.6 | ug/kg | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 360 | 57 | ug/kg | |
| 67-72-1 | Hexachloroethane | ND | 180 | 12 | ug/kg | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 36 | 19 | ug/kg | |
| 78-59-1 | Isophorone | ND | 72 | 6.8 | ug/kg | |
| 91-57-6 | 2-Methylnaphthalene | ND | 72 | 6.8 | ug/kg | |
| 88-74-4 | 2-Nitroaniline | ND | 180 | 8.2 | ug/kg | |
| 99-09-2 | 3-Nitroaniline | ND | 180 | 10 | ug/kg | |
| 100-01-6 | 4-Nitroaniline | ND | 180 | 12 | ug/kg | |
| 91-20-3 | Naphthalene | ND | 36 | 5.8 | ug/kg | |
| 98-95-3 | Nitrobenzene | ND | 72 | 11 | ug/kg | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 72 | 11 | ug/kg | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 180 | 19 | ug/kg | |
| 85-01-8 | Phenanthrene | ND | 36 | 4.0 | ug/kg | |
| 129-00-0 | Pyrene | ND | 36 | 4.5 | ug/kg | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 180 | 8.7 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 367-12-4 | 2-Fluorophenol | 78% | | 30-1 | 06% | |
| | | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



34 of 1690

ACCUTEST JC7512

4.3 4

Page 2 of 3

Report of Analysis

| Client Sample ID: | SB-2-1 | | |
|-------------------|--|---------------|----------|
| Lab Sample ID: | JC7512-3 Da | te Sampled: | 10/29/15 |
| Matrix: | SO - Soil Da | te Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 Pe | rcent Solids: | 87.6 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 4165-62-2 | Phenol-d5 | 75% | | 30-106% |
| 118-79-6 | 2,4,6-Tribromophenol | 89% | | 24-140% |
| 4165-60-0 | Nitrobenzene-d5 | 100% | | 26-122% |
| 321-60-8 | 2-Fluorobiphenyl | 88% | | 36-112% |
| 1718-51-0 | Terphenyl-d14 | 86% | | 36-132% |

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



JC7512

Page 3 of 3

4.3 4

| | Report of Analysis | | | | | | | | | | |
|--|-----------------------------------|--------------------------|------------------------------|------------------|---------------------------|------------------------------|------------------------------------|--|--|--|--|
| Client San Lab Samp Matrix: Method: Project: | le ID: JC7512 SO - So SW846 | oil 8151 SV | W846 3550C ey Avenue, Bro | ooklyn, N | Da Pe | ate Received: 10 | 0/29/15 0/31/15 7.6 | | | | |
| Run #1 Run #2 | File ID OA112010.D | DF 2 | Analyzed 11/05/15 | By VDT | Prep Date 11/03/15 | Prep Batch OP88618 | Analytical Batch GOA3882 | | | | |
| Run #1 Run #2 | Initial Weight 15.2 g | Final V 5.0 ml | <i>T</i> olume | | | | | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|------------------|------------------|-------------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 38 7.5 7.5 | 17 2.2 2.8 | ug/kg ug/kg ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2,4-DCAA | 22% 10% | | 10-1: 10-1: | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| | Report of Analysis Page 1 of | | | | | | | | | | |
|--|---------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|--|-------------------------------|--|--|--|--|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC751 SO - S SW84 | 2-3 oil 5 8081B | SW846 3546 sey Avenue, Bro | ooklyn, N | | Date Sampled: Date Received: Percent Solids: | 10/29/15 10/31/15 87.6 | | | | |
| Run #1 Run #2 | File ID 1G116789.D | DF 1 | Analyzed 11/05/15 | By RK | Prep Date 11/03/15 | Prep Batcl OP88620 | n Analytical Batch G1G3831 | | | | |
| Run #1 Run #2 | Initial Weight 16.1 g | Final V 10.0 m | | | | | | | | | |

Pesticide TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------|--------|--------|------|-------|---|
| 309-00-2 | Aldrin | ND | 0.71 | 0.63 | ug/kg | |
| 319-84-6 | alpha-BHC | ND | 0.71 | 0.47 | ug/kg | |
| 319-85-7 | beta-BHC | ND | 0.71 | 0.44 | ug/kg | |
| 319-86-8 | delta-BHC | ND | 0.71 | 0.28 | ug/kg | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.71 | 0.32 | ug/kg | |
| 5103-71-9 | alpha-Chlordane | ND | 0.71 | 0.38 | ug/kg | |
| 5103-74-2 | gamma-Chlordane | ND | 0.71 | 0.54 | ug/kg | |
| 60-57-1 | Dieldrin | ND | 0.71 | 0.56 | ug/kg | |
| 72-54-8 | 4,4'-DDD | ND | 0.71 | 0.26 | ug/kg | |
| 72-55-9 | 4,4'-DDE | ND | 0.71 | 0.24 | ug/kg | |
| 50-29-3 | 4,4'-DDT | ND | 0.71 | 0.27 | ug/kg | |
| 72-20-8 | Endrin | ND | 0.71 | 0.25 | ug/kg | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.71 | 0.40 | ug/kg | |
| 7421-93-4 | Endrin aldehyde | ND | 0.71 | 0.53 | ug/kg | |
| 959-98-8 | Endosulfan-I | ND | 0.71 | 0.23 | ug/kg | |
| 33213-65-9 | Endosulfan-II | ND | 0.71 | 0.67 | ug/kg | |
| 76-44-8 | Heptachlor | ND | 0.71 | 0.58 | ug/kg | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.71 | 0.29 | ug/kg | |
| 72-43-5 | Methoxychlor | ND | 1.4 | 0.39 | ug/kg | |
| 53494-70-5 | Endrin ketone | ND | 0.71 | 0.37 | ug/kg | |
| 8001-35-2 | Toxaphene | ND | 18 | 12 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 877-09-8 | Tetrachloro-m-xylene | 68% | | 24-1 | 36% | |
| 877-09-8 | Tetrachloro-m-xylene | 68% | | 24-1 | 36% | |
| 2051-24-3 | Decachlorobiphenyl | 43% | | 10-1 | 53% | |
| 2051-24-3 | Decachlorobiphenyl | 82% | | 10-1 | 53% | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

- E = Indicates value exceeds calibration range
- J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



37 of 1690 ACCUTEST JC7512

4.3

| | Report of Analysis Page 1 of 2 | | | | | | | | | | |
|--|---------------------------------------|----------------|-------------------------------|-----------|---------------------------|-----------------------|-------------------------------|--|--|--|--|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC7512 SO - So SW846 | oil 8082A | SW846 3546 sey Avenue, Bro | ooklyn, N | D P | | 10/29/15 10/31/15 87.6 | | | | |
| Run #1 | File ID XX179844.D | DF 1 | Analyzed 11/06/15 | By JR | Prep Date 11/03/15 | Prep Batcl OP88619 | n Analytical Batch GXX5510 | | | | |
| Run #2 | | - | 11,00,10 | | 11,00,10 | 010001/ | | | | | |
| | Initial Weight | | olume | | | | | | | | |
| Run #1 Run #2 | 15.1 g | 10.0 m | 1 | | | | | | | | |

PCB List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|--|--|--|--|--|---|---|
| 12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 11096-82-5 11100-14-4 | Aroclor 1016 Aroclor 1221 Aroclor 1232 Aroclor 1242 Aroclor 1248 Aroclor 1254 Aroclor 1260 Aroclor 1268 | ND ND ND ND ND ND ND | 38 38 38 38 38 38 38 38 38 38 | 12 22 12 17 12 17 16 12 | ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg | |
| 37324-23-5 | Aroclor 1262 | ND | 38 | 11 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 877-09-8 877-09-8 2051-24-3 2051-24-3 | Tetrachloro-m-xylene Tetrachloro-m-xylene Decachlorobiphenyl Decachlorobiphenyl | 93% 93% 85% 86% | | 20-1 20-1 12-1 12-1 | 52% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



38 of 1690 ACCUTEST JC7512

4.3

| Client Sample ID: | SB-2-1 | | |
|-------------------|--|------------------------|----------|
| - | JC7512-3 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| | | Percent Solids: | 87.6 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|---------|-------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | 4950 | 55 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Antimony | < 2.2 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Arsenic | < 2.2 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Barium | < 22 | 22 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C 2 | SW846 3050B ³ |
| Beryllium | 0.28 | 0.22 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Cadmium | < 0.55 | 0.55 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Calcium | 1460 | 550 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Chromium | 13.8 | 1.1 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Cobalt | 32.3 | 5.5 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Copper | 8.1 | 2.8 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Iron | 10400 | 55 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Lead | 4.4 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Magnesium | 2780 | 550 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Manganese | 210 | 1.7 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Mercury | < 0.035 | 0.035 | mg/kg | 1 | 11/04/15 | 11/04/15 ма | SW846 7471B ¹ | SW846 7471B ⁴ |
| Nickel | 37.0 | 4.4 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Potassium | < 1100 | 1100 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Selenium | < 2.2 | 2.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Silver | 1.0 | 0.55 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Sodium | < 1100 | 1100 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Thallium | < 1.1 | 1.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Vanadium | 15.2 | 5.5 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Zinc | 19.5 | 5.5 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |

(1) Instrument QC Batch: MA37945

(2) Instrument QC Batch: MA37978

(3) Prep QC Batch: MP90046

(4) Prep QC Batch: MP90074





JC7512

| | | | 1 | | v | | e |
|------------------|-------------|--------------|-----------------|------------|-----------|-------------------|------------------|
| Client Sa | mple ID: SE | 8-2-5 | | | | | |
| Lab Sam | ple ID: JC | 7512-4 | | | Da | ate Sampled: 10 |)/29/15 |
| Matrix: | SC |) - Soil | | | Da | ate Received: 10 |)/31/15 |
| Method: | SV | V846 8260C | | | Pe | ercent Solids: 96 | 5.6 |
| Project: | 20 | 02-2024 Crop | sey Avenue, Bro | ooklyn, NY | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 Run #2 | I209242.D | 1 | 11/03/15 | SĴM | n/a | n/a | VI8396 |
| | Initial Wei | ght | | | | | |
| Run #1 Run #2 | 4.9 g | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 19.6 | 11 | 2.4 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.53 | 0.14 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.3 | 0.33 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.1 | 0.16 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.3 | 0.25 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.3 | 0.38 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 11 | 2.0 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.1 | 0.24 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.1 | 0.24 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.1 | 0.16 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.3 | 0.51 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.1 | 0.16 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.3 | 0.28 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.1 | 0.33 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.1 | 0.57 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.1 | 0.22 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.1 | 0.14 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.1 | 0.13 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.1 | 0.17 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.1 | 0.24 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.3 | 0.38 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.1 | 0.15 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.1 | 0.14 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.1 | 0.63 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.1 | 0.82 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.1 | 0.63 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.1 | 0.25 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.1 | 0.12 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.1 | 0.19 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.1 | 0.17 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.3 | 0.47 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.3 | 1.4 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



JC7512

4.4 **4**

Page 1 of 2

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Samp Lab Sample Matrix: Method: Project: | | Avenue, Brool | klyn, NY | | Date | Sampled: Received: ent Solids: | 10/29/15 10/31/15 96.6 | |
|---|---------------------------|---------------|----------|------|-------|--------------------------------------|------------------------------|--|
| VOA TCL I | List | | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | | |
| 98-82-8 | Isopropylbenzene | ND | 2.1 | 0.11 | ug/kg | | | |
| 79-20-9 | Methyl Acetate | ND | 5.3 | 0.91 | ug/kg | | | |
| 108-87-2 | Methylcyclohexane | ND | 2.1 | 0.24 | ug/kg | | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.1 | 0.16 | ug/kg | | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIB | K) ND | 5.3 | 0.48 | ug/kg | | | |
| 75-09-2 | Methylene chloride | ND | 5.3 | 1.0 | ug/kg | | | |
| 100-42-5 | Styrene | ND | 2.1 | 0.19 | ug/kg | | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.1 | 0.18 | ug/kg | | | |
| 127-18-4 | Tetrachloroethene | 44.8 | 2.1 | 0.32 | ug/kg | | | |
| 108-88-3 | Toluene | ND | 1.1 | 0.22 | ug/kg | | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.3 | 0.19 | ug/kg | | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.3 | 0.18 | ug/kg | | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.1 | 0.16 | ug/kg | | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.1 | 0.16 | ug/kg | | | |
| 79-01-6 | Trichloroethene | 0.88 | 1.1 | 0.16 | ug/kg | J | | |
| 75-69-4 | Trichlorofluoromethane | ND | 5.3 | 0.26 | ug/kg | | | |
| 75-01-4 | Vinyl chloride | ND | 2.1 | 0.21 | ug/kg | | | |
| | m,p-Xylene | ND | 1.1 | 0.37 | ug/kg | | | |
| 95-47-6 | o-Xylene | ND | 1.1 | 0.29 | ug/kg | | | |
| 1330-20-7 | Xylene (total) | ND | 1.1 | 0.29 | ug/kg | | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | | |
| 1868-53-7 | Dibromofluoromethane | 101% | | 70-1 | 22% | | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 98% | | | 24% | | | |
| 2037-26-5 | Toluene-D8 | 92% | | | 25% | | | |
| | | | | | | | | |

94%

Page 2 of 2

4.4

4-Bromofluorobenzene

460-00-4

J = Indicates an estimated value

72-130%

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



41 of 1690 ACCUTEST JC7512

| | Report of Analysis Pa | | | | | | | | | | |
|---|---------------------------------|----------------------------|-------------------------------|-----------------|---------------------------|-----------------------|------------------------------------|--|--|--|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC751 SO - S SW84 | 12-4 Soil 16 8270D | SW846 3546 sey Avenue, Bro | ooklyn, N | Da Pe | te Received: 10 |)/29/15)/31/15 5.6 | | | | |
| Run #1 Run #2 | File ID 3E77586.D | DF 1 | Analyzed 11/03/15 | By AN | Prep Date 11/03/15 | Prep Batch OP88613 | Analytical Batch E3E3376 | | | | |
| Run #1 Run #2 | Initial Weigh 30.9 g | t Final V 1.0 ml | olume | | | | | | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units Q |) |
|-----------|----------------------------|--------|-----|-----|---------|---|
| 95-57-8 | 2-Chlorophenol | ND | 67 | 25 | ug/kg | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 170 | 30 | ug/kg | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 170 | 27 | ug/kg | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 170 | 61 | ug/kg | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 170 | 150 | ug/kg | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 170 | 64 | ug/kg | |
| 95-48-7 | 2-Methylphenol | ND | 67 | 49 | ug/kg | |
| | 3&4-Methylphenol | ND | 67 | 32 | ug/kg | |
| 88-75-5 | 2-Nitrophenol | ND | 170 | 31 | ug/kg | |
| 100-02-7 | 4-Nitrophenol | ND | 340 | 57 | ug/kg | |
| 87-86-5 | Pentachlorophenol | ND | 170 | 82 | ug/kg | |
| 108-95-2 | Phenol | ND | 67 | 25 | ug/kg | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 170 | 32 | ug/kg | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 170 | 30 | ug/kg | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 170 | 27 | ug/kg | |
| 83-32-9 | Acenaphthene | ND | 34 | 32 | ug/kg | |
| 208-96-8 | Acenaphthylene | ND | 34 | 3.5 | ug/kg | |
| 98-86-2 | Acetophenone | ND | 170 | 5.7 | ug/kg | |
| 120-12-7 | Anthracene | ND | 34 | 2.9 | ug/kg | |
| 1912-24-9 | Atrazine | ND | 67 | 14 | ug/kg | |
| 56-55-3 | Benzo(a)anthracene | ND | 34 | 6.5 | ug/kg | |
| 50-32-8 | Benzo(a)pyrene | ND | 34 | 7.1 | ug/kg | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 34 | 6.9 | ug/kg | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 34 | 10 | ug/kg | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 34 | 7.5 | ug/kg | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 67 | 7.6 | ug/kg | |
| 85-68-7 | Butyl benzyl phthalate | ND | 67 | 18 | ug/kg | |
| 92-52-4 | 1,1'-Biphenyl | ND | 67 | 6.2 | ug/kg | |
| 100-52-7 | Benzaldehyde | ND | 170 | 8.4 | ug/kg | |
| 91-58-7 | 2-Chloronaphthalene | ND | 67 | 4.8 | ug/kg | |
| 106-47-8 | 4-Chloroaniline | ND | 170 | 8.9 | ug/kg | |
| 86-74-8 | Carbazole | ND | 67 | 3.7 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



4.4



E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Sample ID: | SB-2-5 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-4 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 96.6 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|--------|------|-------|---|
| 105-60-2 | Caprolactam | ND | 67 | 22 | ug/kg | |
| 218-01-9 | Chrysene | ND | 34 | 5.4 | ug/kg | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 67 | 7.6 | ug/kg | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 67 | 14 | ug/kg | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 67 | 7.7 | ug/kg | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 67 | 6.3 | ug/kg | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 34 | 6.3 | ug/kg | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 34 | 8.6 | ug/kg | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 67 | 22 | ug/kg | |
| 123-91-1 | 1,4-Dioxane | ND | 34 | 23 | ug/kg | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 34 | 12 | ug/kg | |
| 132-64-9 | Dibenzofuran | ND | 67 | 4.7 | ug/kg | |
| 84-74-2 | Di-n-butyl phthalate | ND | 67 | 4.0 | ug/kg | |
| 117-84-0 | Di-n-octyl phthalate | ND | 67 | 4.5 | ug/kg | |
| 84-66-2 | Diethyl phthalate | ND | 67 | 4.3 | ug/kg | |
| 131-11-3 | Dimethyl phthalate | ND | 67 | 4.8 | ug/kg | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 67 | 12 | ug/kg | |
| 206-44-0 | Fluoranthene | ND | 34 | 4.1 | ug/kg | |
| 86-73-7 | Fluorene | ND | 34 | 4.0 | ug/kg | |
| 118-74-1 | Hexachlorobenzene | ND | 67 | 6.6 | ug/kg | |
| 87-68-3 | Hexachlorobutadiene | ND | 34 | 8.9 | ug/kg | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 340 | 53 | ug/kg | |
| 67-72-1 | Hexachloroethane | ND | 170 | 11 | ug/kg | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 34 | 17 | ug/kg | |
| 78-59-1 | Isophorone | ND | 67 | 6.3 | ug/kg | |
| 91-57-6 | 2-Methylnaphthalene | ND | 67 | 6.3 | ug/kg | |
| 88-74-4 | 2-Nitroaniline | ND | 170 | 7.6 | ug/kg | |
| 99-09-2 | 3-Nitroaniline | ND | 170 | 9.5 | ug/kg | |
| 100-01-6 | 4-Nitroaniline | ND | 170 | 11 | ug/kg | |
| 91-20-3 | Naphthalene | ND | 34 | 5.4 | ug/kg | |
| 98-95-3 | Nitrobenzene | ND | 67 | 11 | ug/kg | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 67 | 9.9 | ug/kg | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 170 | 18 | ug/kg | |
| 85-01-8 | Phenanthrene | ND | 34 | 3.7 | ug/kg | |
| 129-00-0 | Pyrene | ND | 34 | 4.2 | ug/kg | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 170 | 8.0 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 367-12-4 | 2-Fluorophenol | 79% | | 30-1 | 06% | |
| | | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound

Page 2 of 3



43 of 1690

ACCUTEST JC7512

Report of Analysis

| Client Sample ID: | SB-2-5 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-4 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 96.6 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 4165-62-2 | Phenol-d5 | 78% | | 30-106% |
| 118-79-6 | 2,4,6-Tribromophenol | 92% | | 24-140% |
| 4165-60-0 | Nitrobenzene-d5 | 101% | | 26-122% |
| 321-60-8 | 2-Fluorobiphenyl | 90% | | 36-112% |
| 1718-51-0 | Terphenyl-d14 | 85% | | 36-132% |

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 3 of 3



| | Report of Analysis | | | | | | | | |
|--|------------------------------------|------------------------|--------------------------------|------------------|---------------------------|--|------------------------------|--|--|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: JC7512 SO - So SW846 | oil 8151 S | W846 3550C osey Avenue, Bro | ooklyn, NY |] | Date Sampled: Date Received: Percent Solids: | 10/29/15 10/31/15 96.6 | | |
| Run #1 Run #2 | File ID OA112011.D | DF 2 | Analyzed 11/05/15 | By VDT | Prep Date 11/03/15 | Prep Batcl OP88618 | Analytical Batch GOA3882 | | |
| Run #1 Run #2 | Initial Weight 15.0 g | Final 5.0 ml | Volume | | | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|------------------|------------------|-------------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 35 6.9 6.9 | 15 2.0 2.6 | ug/kg ug/kg ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2.4-DCAA | 23% 79% | | 10-1 10-1 | | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit E = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| | Report of Analysis Pa | | | | | | | | |
|--|-------------------------------|------------------------------|--------------------------------|-----------------|---------------------------|-----------------------|-----------------------------|--|--|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: JC7 SO SW | 512-4 - Soil 846 8081B | SW846 3546 psey Avenue, Bro | ooklyn, 1 | Da Pe | te Received: 10 | 0/29/15 0/31/15 5.6 | | |
| Run #1 Run #2 | File ID 1G116790.I | DF 0 1 | Analyzed 11/05/15 | By RK | Prep Date 11/03/15 | Prep Batch OP88620 | Analytical Batch G1G3831 | | |
| Run #1 Run #2 | Initial Weig 16.5 g | g ht Final 10.0 1 | Volume nl | | | | | | |

Pesticide TCL List

| CAS No. Compound | | Result | RL | MDL | Units | Q |
|------------------|----------------------|--------|--------|------|-------|---|
| 309-00-2 | Aldrin | ND | 0.63 | 0.56 | ug/kg | |
| 319-84-6 | alpha-BHC | ND | 0.63 | 0.42 | ug/kg | |
| 319-85-7 | beta-BHC | ND | 0.63 | 0.39 | ug/kg | |
| 319-86-8 | delta-BHC | ND | 0.63 | 0.25 | ug/kg | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.63 | 0.29 | ug/kg | |
| 5103-71-9 | alpha-Chlordane | ND | 0.63 | 0.33 | ug/kg | |
| 5103-74-2 | gamma-Chlordane | ND | 0.63 | 0.48 | ug/kg | |
| 60-57-1 | Dieldrin | ND | 0.63 | 0.49 | ug/kg | |
| 72-54-8 | 4,4'-DDD | ND | 0.63 | 0.23 | ug/kg | |
| 72-55-9 | 4,4'-DDE | ND | 0.63 | 0.21 | ug/kg | |
| 50-29-3 | 4,4'-DDT | ND | 0.63 | 0.24 | ug/kg | |
| 72-20-8 | Endrin | ND | 0.63 | 0.22 | ug/kg | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.63 | 0.36 | ug/kg | |
| 7421-93-4 | Endrin aldehyde | ND | 0.63 | 0.47 | ug/kg | |
| 959-98-8 | Endosulfan-I | ND | 0.63 | 0.21 | ug/kg | |
| 33213-65-9 | Endosulfan-II | ND | 0.63 | 0.59 | ug/kg | |
| 76-44-8 | Heptachlor | ND | 0.63 | 0.52 | ug/kg | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.63 | 0.26 | ug/kg | |
| 72-43-5 | Methoxychlor | ND | 1.3 | 0.35 | ug/kg | |
| 53494-70-5 | Endrin ketone | ND | 0.63 | 0.33 | ug/kg | |
| 8001-35-2 | Toxaphene | ND | 16 | 11 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 877-09-8 | Tetrachloro-m-xylene | 85% | | 24-1 | 36% | |
| 877-09-8 | Tetrachloro-m-xylene | 87% | | 24-1 | 36% | |
| 2051-24-3 | Decachlorobiphenyl | 59% | | 10-1 | 53% | |
| 2051-24-3 | Decachlorobiphenyl | 100% | | 10-1 | 53% | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

- E = Indicates value exceeds calibration range
- J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



46 of 1690 ACCUTEST JC7512

| | Report of Analysis P | | | | | | | | |
|--|---------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|-----------------------|------------------------------|--|--|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC751 SO - S SW84 | 2-4 oil 6 8082A | SW846 3546 sey Avenue, Bro | ooklyn, N | D P | ate Received: | 10/29/15 10/31/15 96.6 | | |
| Run #1 Run #2 | File ID XX179849.D | DF 1 | Analyzed 11/07/15 | By JR | Prep Date 11/03/15 | Prep Batch OP88619 | Analytical Batch GXX5510 | | |
| Run #1 Run #2 | Initial Weight 16.5 g | Final V 10.0 m | | | | | | | |

PCB List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------|--------|--------|------|-------|---|
| 12674-11-2 | Aroclor 1016 | ND | 31 | 10 | ug/kg | |
| 11104-28-2 | Aroclor 1221 | ND | 31 | 18 | ug/kg | |
| 11141-16-5 | Aroclor 1232 | ND | 31 | 10 | ug/kg | |
| 53469-21-9 | Aroclor 1242 | ND | 31 | 14 | ug/kg | |
| 12672-29-6 | Aroclor 1248 | ND | 31 | 9.6 | ug/kg | |
| 11097-69-1 | Aroclor 1254 | ND | 31 | 14 | ug/kg | |
| 11096-82-5 | Aroclor 1260 | ND | 31 | 13 | ug/kg | |
| 11100-14-4 | Aroclor 1268 | ND | 31 | 9.7 | ug/kg | |
| 37324-23-5 | Aroclor 1262 | ND | 31 | 8.8 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| | 8 | | | | | |
| 877-09-8 | Tetrachloro-m-xylene | 110% | | 20-1 | 52% | |
| 877-09-8 | Tetrachloro-m-xylene | 112% | | 20-1 | 52% | |
| 2051-24-3 | Decachlorobiphenyl | 98% | | 12-1 | 57% | |
| 2051-24-3 | Decachlorobiphenyl | 101% | | 12-1 | 57% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



47 of 1690 ACCUTEST JC7512

4.4

| Client Sample ID: | SB-2-5 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-4 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| | | Percent Solids: | 96.6 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|---------|-------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | 4200 | 50 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Antimony | < 2.0 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Arsenic | < 2.0 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Barium | 21.6 | 20 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Beryllium | 0.31 | 0.20 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Cadmium | < 0.50 | 0.50 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Calcium | 868 | 500 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Chromium | 13.9 | 1.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Cobalt | 5.8 | 5.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Copper | 8.7 | 2.5 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Iron | 9930 | 50 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Lead | 3.8 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Magnesium | 2090 | 500 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Manganese | 239 | 1.5 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Mercury | < 0.033 | 0.033 | mg/kg | 1 | 11/04/15 | 11/04/15 ма | SW846 7471B ¹ | SW846 7471B ⁴ |
| Nickel | 41.5 | 4.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Potassium | < 1000 | 1000 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Selenium | < 2.0 | 2.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Silver | < 0.50 | 0.50 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Sodium | < 1000 | 1000 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Thallium | < 1.0 | 1.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Vanadium | 14.0 | 5.0 | mg/kg | 1 | 11/03/15 | 11/07/15 MS | SW846 6010C ² | SW846 3050B ³ |
| Zinc | 19.1 | 5.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |

(1) Instrument QC Batch: MA37945

(2) Instrument QC Batch: MA37978

(3) Prep QC Batch: MP90046

(4) Prep QC Batch: MP90074



Page 1 of 1

48 of 1690

ACCUTEST JC7512

| | | | Repo | ort of A | nalysis | | Page 1 of 2 |
|--|-----------------------------|--------------------------|-----------------------------|------------------|-------------------------|--|------------------------------|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC75 SO - SW8 | 12-5 Soil 46 8260C | sey Avenue, Bro | ooklyn, NY | | Date Sampled: Date Received: Percent Solids: | 10/29/15 10/31/15 89.5 |
| Run #1 Run #2 | File ID I209243.D | DF 1 | Analyzed 11/03/15 | By SJM | Prep Date n/a | Prep Batc n/a | h Analytical Batch V18396 |
| Run #1 Run #2 | Initial Weigh 5.2 g | ıt | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 14.0 | 11 | 2.4 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.54 | 0.14 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.4 | 0.33 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.1 | 0.17 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.4 | 0.25 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.4 | 0.39 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 11 | 2.1 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.1 | 0.24 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.1 | 0.25 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.1 | 0.17 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.4 | 0.52 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.1 | 0.16 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.4 | 0.28 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.1 | 0.34 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.1 | 0.58 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.1 | 0.22 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.1 | 0.14 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.1 | 0.13 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.1 | 0.17 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.1 | 0.24 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.4 | 0.39 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.1 | 0.15 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.1 | 0.14 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.1 | 0.64 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.1 | 0.84 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.1 | 0.64 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.1 | 0.26 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.1 | 0.13 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.1 | 0.19 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.1 | 0.18 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.4 | 0.48 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.4 | 1.4 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



49 of 1690 SGS ACCUTEST

JC7512

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Sample ID: Lab Sample ID: | JC7512-5 | Date Sampled: | |
|-------------------------------------|--|-----------------|----------|
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8260C | Percent Solids: | 89.5 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 2.1 | 0.11 | ug/kg | |
| 79-20-9 | Methyl Acetate | ND | 5.4 | 0.92 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 2.1 | 0.24 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.1 | 0.16 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.4 | 0.49 | ug/kg | |
| 75-09-2 | Methylene chloride | ND | 5.4 | 1.1 | ug/kg | |
| 100-42-5 | Styrene | ND | 2.1 | 0.19 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.1 | 0.19 | ug/kg | |
| 127-18-4 | Tetrachloroethene | 23.3 | 2.1 | 0.32 | ug/kg | |
| 108-88-3 | Toluene | ND | 1.1 | 0.22 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.4 | 0.19 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.4 | 0.18 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.1 | 0.16 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.1 | 0.16 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 1.1 | 0.16 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 5.4 | 0.27 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 2.1 | 0.21 | ug/kg | |
| | m,p-Xylene | ND | 1.1 | 0.38 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 1.1 | 0.29 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 1.1 | 0.29 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 99% | | 70-12 | 22% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 97% | | 68-12 | 24% | |
| 2037-26-5 | Toluene-D8 | 94% | | 77-12 | 25% | |
| 460-00-4 | 4-Bromofluorobenzene | 91% | | 72-13 | 30% | |

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



50 of 1690

ACCUTEST JC7512

Page 2 of 2

4.5

4

| | Report of Analysis | | | | | | | | |
|---|-----------------------------------|-----------------------|------------------------------|-----------|---------------------------|-----------------------|------------------------------------|--|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC7512 SO - S SW846 | 2-5 oil 5 8270D | SW846 3546 ey Avenue, Bro | ooklyn, N | Da Pe | · · · · · · |)/29/15)/31/15).5 | | |
| Run #1 Run #2 | File ID 3E77587.D | DF 1 | Analyzed 11/03/15 | By An | Prep Date 11/03/15 | Prep Batch OP88613 | Analytical Batch E3E3376 | | |
| Run #1 Run #2 | Initial Weight 30.9 g | Final V 1.0 ml | olume | | | | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|----------------------------|--------|-----|-----|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 72 | 27 | ug/kg | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 180 | 33 | ug/kg | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 180 | 29 | ug/kg | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 180 | 66 | ug/kg | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 180 | 160 | ug/kg | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 180 | 69 | ug/kg | |
| 95-48-7 | 2-Methylphenol | ND | 72 | 52 | ug/kg | |
| | 3&4-Methylphenol | ND | 72 | 35 | ug/kg | |
| 88-75-5 | 2-Nitrophenol | ND | 180 | 33 | ug/kg | |
| 100-02-7 | 4-Nitrophenol | ND | 360 | 61 | ug/kg | |
| 87-86-5 | Pentachlorophenol | ND | 180 | 88 | ug/kg | |
| 108-95-2 | Phenol | ND | 72 | 27 | ug/kg | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 180 | 34 | ug/kg | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 180 | 33 | ug/kg | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 180 | 29 | ug/kg | |
| 83-32-9 | Acenaphthene | ND | 36 | 34 | ug/kg | |
| 208-96-8 | Acenaphthylene | ND | 36 | 3.8 | ug/kg | |
| 98-86-2 | Acetophenone | ND | 180 | 6.1 | ug/kg | |
| 120-12-7 | Anthracene | ND | 36 | 3.1 | ug/kg | |
| 1912-24-9 | Atrazine | ND | 72 | 15 | ug/kg | |
| 56-55-3 | Benzo(a)anthracene | ND | 36 | 7.0 | ug/kg | |
| 50-32-8 | Benzo(a)pyrene | ND | 36 | 7.7 | ug/kg | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 36 | 7.4 | ug/kg | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 36 | 11 | ug/kg | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 36 | 8.1 | ug/kg | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 72 | 8.2 | ug/kg | |
| 85-68-7 | Butyl benzyl phthalate | ND | 72 | 20 | ug/kg | |
| 92-52-4 | 1,1'-Biphenyl | ND | 72 | 6.7 | ug/kg | |
| 100-52-7 | Benzaldehyde | ND | 180 | 9.0 | ug/kg | |
| 91-58-7 | 2-Chloronaphthalene | ND | 72 | 5.2 | ug/kg | |
| 106-47-8 | 4-Chloroaniline | ND | 180 | 9.6 | ug/kg | |
| 86-74-8 | Carbazole | ND | 72 | 4.0 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



| Client Sample ID: | SB-3-1 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-5 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 89.5 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|--------|------|-------|---|
| 105-60-2 | Caprolactam | ND | 72 | 23 | ug/kg | |
| 218-01-9 | Chrysene | ND | 36 | 5.8 | ug/kg | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 72 | 8.2 | ug/kg | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 72 | 15 | ug/kg | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 72 | 8.3 | ug/kg | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 72 | 6.8 | ug/kg | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 36 | 6.8 | ug/kg | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 36 | 9.3 | ug/kg | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 72 | 24 | ug/kg | |
| 123-91-1 | 1,4-Dioxane | ND | 36 | 24 | ug/kg | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 36 | 13 | ug/kg | |
| 132-64-9 | Dibenzofuran | ND | 72 | 5.0 | ug/kg | |
| 84-74-2 | Di-n-butyl phthalate | ND | 72 | 4.3 | ug/kg | |
| 117-84-0 | Di-n-octyl phthalate | ND | 72 | 4.9 | ug/kg | |
| 84-66-2 | Diethyl phthalate | ND | 72 | 4.6 | ug/kg | |
| 131-11-3 | Dimethyl phthalate | ND | 72 | 5.2 | ug/kg | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 72 | 13 | ug/kg | |
| 206-44-0 | Fluoranthene | ND | 36 | 4.4 | ug/kg | |
| 86-73-7 | Fluorene | ND | 36 | 4.3 | ug/kg | |
| 118-74-1 | Hexachlorobenzene | ND | 72 | 7.1 | ug/kg | |
| 87-68-3 | Hexachlorobutadiene | ND | 36 | 9.6 | ug/kg | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 360 | 57 | ug/kg | |
| 67-72-1 | Hexachloroethane | ND | 180 | 12 | ug/kg | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 36 | 19 | ug/kg | |
| 78-59-1 | Isophorone | ND | 72 | 6.8 | ug/kg | |
| 91-57-6 | 2-Methylnaphthalene | ND | 72 | 6.8 | ug/kg | |
| 88-74-4 | 2-Nitroaniline | ND | 180 | 8.2 | ug/kg | |
| 99-09-2 | 3-Nitroaniline | ND | 180 | 10 | ug/kg | |
| 100-01-6 | 4-Nitroaniline | ND | 180 | 12 | ug/kg | |
| 91-20-3 | Naphthalene | ND | 36 | 5.8 | ug/kg | |
| 98-95-3 | Nitrobenzene | ND | 72 | 11 | ug/kg | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 72 | 11 | ug/kg | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 180 | 19 | ug/kg | |
| 85-01-8 | Phenanthrene | ND | 36 | 4.0 | ug/kg | |
| 129-00-0 | Pyrene | ND | 36 | 4.5 | ug/kg | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 180 | 8.7 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 367-12-4 | 2-Fluorophenol | 83% | | 30-1 | 06% | |
| | | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound





Report of Analysis

| Client Sample ID: | SB-3-1 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC7512-5 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 89.5 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 4165-62-2 | Phenol-d5 | 82% | | 30-106% |
| 118-79-6 | 2,4,6-Tribromophenol | 98% | | 24-140% |
| 4165-60-0 | Nitrobenzene-d5 | 109% | | 26-122% |
| 321-60-8 | 2-Fluorobiphenyl | 93% | | 36-112% |
| 1718-51-0 | Terphenyl-d14 | 91% | | 36-132% |

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



SGS

4.5

Page 3 of 3

ACCUTEST JC7512

| | Report of Analysis | | | | | | | |
|--|------------------------------------|-------------------|------------------------------|------------------|---------------------------|------------------------------|------------------------------------|--|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: JC7512 SO - So SW846 | il 8151 SW | V846 3550C ey Avenue, Bro | oklyn, NY | Da Pe | ate Received: 10 |)/29/15)/31/15 9.5 | |
| Run #1 Run #2 | File ID OA112012.D | DF 2 | Analyzed 11/05/15 | By VDT | Prep Date 11/03/15 | Prep Batch OP88618 | Analytical Batch GOA3882 | |
| Run #1 Run #2 | Initial Weight 15.8 g | Final V 5.0 ml | olume | | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|------------------|------------------|-------------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 35 7.1 7.1 | 16 2.1 2.6 | ug/kg ug/kg ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2,4-DCAA | 19% 16% | | 10-1 10-1 | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| Report of Analysis Page 1 | | | | | | | | | | |
|--|---------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|------------------------------|------------------------------------|--|--|--|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC751 SO - S SW84 | 2-5 oil 6 8081B | SW846 3546 sey Avenue, Bro | ooklyn, N | I F | Date Received: | 10/29/15 10/31/15 89.5 | | | |
| Run #1 Run #2 | File ID 1G116791.D | DF 1 | Analyzed 11/05/15 | By RK | Prep Date 11/03/15 | Prep Batch OP88620 | Analytical Batch G1G3831 | | | |
| Run #1 Run #2 | Initial Weight 15.2 g | Final V 10.0 m | | | | | | | | |

Pesticide TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------|--------|---------|------|-------|---|
| 309-00-2 | Aldrin | ND | 0.74 | 0.66 | ug/kg | |
| 319-84-6 | alpha-BHC | ND | 0.74 | 0.49 | ug/kg | |
| 319-85-7 | beta-BHC | ND | 0.74 | 0.45 | ug/kg | |
| 319-86-8 | delta-BHC | ND | 0.74 | 0.29 | ug/kg | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.74 | 0.33 | ug/kg | |
| 5103-71-9 | alpha-Chlordane | ND | 0.74 | 0.39 | ug/kg | |
| 5103-74-2 | gamma-Chlordane | ND | 0.74 | 0.56 | ug/kg | |
| 60-57-1 | Dieldrin | ND | 0.74 | 0.58 | ug/kg | |
| 72-54-8 | 4,4'-DDD | ND | 0.74 | 0.27 | ug/kg | |
| 72-55-9 | 4,4'-DDE | ND | 0.74 | 0.24 | ug/kg | |
| 50-29-3 | 4,4'-DDT | ND | 0.74 | 0.28 | ug/kg | |
| 72-20-8 | Endrin | ND | 0.74 | 0.26 | ug/kg | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.74 | 0.42 | ug/kg | |
| 7421-93-4 | Endrin aldehyde | ND | 0.74 | 0.55 | ug/kg | |
| 959-98-8 | Endosulfan-I | ND | 0.74 | 0.24 | ug/kg | |
| 33213-65-9 | Endosulfan-II | ND | 0.74 | 0.70 | ug/kg | |
| 76-44-8 | Heptachlor | ND | 0.74 | 0.60 | ug/kg | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.74 | 0.30 | ug/kg | |
| 72-43-5 | Methoxychlor | ND | 1.5 | 0.41 | ug/kg | |
| 53494-70-5 | Endrin ketone | ND | 0.74 | 0.39 | ug/kg | |
| 8001-35-2 | Toxaphene | ND | 18 | 13 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 877-09-8 | Tetrachloro-m-xylene | 79% | | 24-1 | 36% | |
| 877-09-8 | Tetrachloro-m-xylene | 79% | | 24-1 | 36% | |
| 2051-24-3 | Decachlorobiphenyl | 64% | | 10-1 | 53% | |
| 2051-24-3 | Decachlorobiphenyl | 95% | 10-153% | | | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

- E = Indicates value exceeds calibration range
- J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





| Report of Analysis Page 1 | | | | | | | | | | |
|--|------------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|--|-------------------------------|--|--|--|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC7512 SO - So SW846 | oil 8082A | SW846 3546 sey Avenue, Bro | ooklyn, N | I F | Date Sampled: Date Received: Percent Solids: | 10/29/15 10/31/15 89.5 | | | |
| Run #1 | File ID XX179850.D | DF 1 | Analyzed 11/07/15 | By JR | Prep Date 11/03/15 | Prep Batcl OP88619 | n Analytical Batch GXX5510 | | | |
| Run #2 | | | | | | | | | | |
| Run #1 Run #2 | Initial Weight 15.2 g | Final V 10.0 m | | | | | | | | |

PCB List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|--|--|--|--|--|---|---|
| 12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 11096-82-5 11100-14-4 | Aroclor 1016 Aroclor 1221 Aroclor 1232 Aroclor 1242 Aroclor 1248 Aroclor 1254 Aroclor 1260 Aroclor 1268 | ND ND ND ND ND ND ND | 37 37 37 37 37 37 37 37 37 | 12 22 12 17 11 16 16 11 | ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg | |
| 37324-23-5 CAS No. | Aroclor 1262 Surrogate Recoveries | ND Run# 1 | 37 Run# 2 | 10 Lim | | |
| 877-09-8 877-09-8 2051-24-3 2051-24-3 | Tetrachloro-m-xylene Tetrachloro-m-xylene Decachlorobiphenyl Decachlorobiphenyl | 103% 102% 95% 97% | 20-152% 20-152% 12-157% 12-157% | | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| Client Sample ID: | SB-3-1 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC7512-5 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| | | Percent Solids: | 89.5 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|---------|-------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | 4880 | 57 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Antimony | < 2.3 | 2.3 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Arsenic | < 2.3 | 2.3 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Barium | 23.5 | 23 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C 2 | SW846 3050B ³ |
| Beryllium | 0.32 | 0.23 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Cadmium | < 0.57 | 0.57 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Calcium | 1000 | 570 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Chromium | 15.5 | 1.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Cobalt | 8.4 | 5.7 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Copper | 9.5 | 2.9 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Iron | 11100 | 57 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C 2 | SW846 3050B ³ |
| Lead | 5.0 | 2.3 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Magnesium | 2120 | 570 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Manganese | 257 | 1.7 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Mercury | < 0.036 | 0.036 | mg/kg | 1 | 11/04/15 | 11/04/15 ма | SW846 7471B ¹ | SW846 7471B ⁴ |
| Nickel | 43.0 | 4.6 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Potassium | < 1100 | 1100 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Selenium | < 2.3 | 2.3 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Silver | < 0.57 | 0.57 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Sodium | < 1100 | 1100 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Thallium | < 1.1 | 1.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Vanadium | 15.9 | 5.7 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Zinc | 20.8 | 5.7 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |

(1) Instrument QC Batch: MA37945

(2) Instrument QC Batch: MA37978

(3) Prep QC Batch: MP90046

(4) Prep QC Batch: MP90074

4.5 **4**



JC7512

| e ID: SB-3- D: JC751 | - | | | | | |
|-------------------------|---|---|---|--|--|---|
| D: JC751 | 20 | | | | | |
| | 2-0 | | | Da | te Sampled: 10 | 0/29/15 |
| SO - 5 | Soil | | | Da | te Received: 10 | 0/31/15 |
| SW84 | 6 8260C | | | Ре | rcent Solids: 95 | 5.5 |
| 2002- | 2024 Crops | sey Avenue, Bro | ooklyn, NY | | | |
| ile ID | DF | Analyzed | Bv | Prep Date | Prep Batch | Analytical Batch |
| 209245.D | 1 | 11/03/15 | SJM | n/a | n/a | VI8396 |
| | | | | | | |
| nitial Weight | t | | | | | |
| .1 g | | | | | | |
| e | | | | | | |
| | SW84 2002- `ile ID 209245.D | Sile ID DF 209245.D 1 nitial Weight | SW846 8260C 2002-2024 Cropsey Avenue, Browne, Bro | SW846 8260C 2002-2024 Cropsey Avenue, Brooklyn, NY Sile ID DF Analyzed By 209245.D 1 11/03/15 SJM | SW846 8260C Pe 2002-2024 Cropsey Avenue, Brooklyn, NY File ID DF Analyzed By Prep Date 209245.D 1 11/03/15 SJM n/a nitial Weight | SW846 8260C Percent Solids: 9: 2002-2024 Cropsey Avenue, Brooklyn, NY Prep Date DF Analyzed By Prep Date Prep Batch 209245.D 1 11/03/15 SJM n/a n/a |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 11.6 | 10 | 2.3 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.51 | 0.14 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.1 | 0.32 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.1 | 0.16 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.1 | 0.24 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.1 | 0.37 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 2.0 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.1 | 0.23 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.1 | 0.24 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.1 | 0.16 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.1 | 0.49 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.1 | 0.15 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.1 | 0.27 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.1 | 0.32 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.1 | 0.56 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.1 | 0.21 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.13 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.13 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.16 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.23 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.1 | 0.37 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.14 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.14 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.61 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.80 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.61 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.1 | 0.24 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.1 | 0.12 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.1 | 0.18 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.17 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.1 | 0.46 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.1 | 1.4 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2



E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Samp Lab Sample Matrix: Method: Project: | | venue, Broo | klyn, NY | | Date | Sampled: Received: ent Solids: | 10/29/15 10/31/15 95.5 |
|---|----------------------------|-------------|----------|------|-------|--------------------------------------|------------------------------|
| VOA TCL I | | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 2.1 | 0.11 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 5.1 | 0.88 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 2.1 | 0.23 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.16 | ug/kg | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.1 | 0.47 | ug/kg | | |
| 75-09-2 | Methylene chloride | ND | 5.1 | 1.0 | ug/kg | | |
| 100-42-5 | Styrene | ND | 2.1 | 0.18 | ug/kg | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.1 | 0.18 | ug/kg | | |
| 127-18-4 | Tetrachloroethene | 23.8 | 2.1 | 0.31 | ug/kg | | |
| 108-88-3 | Toluene | ND | 1.0 | 0.21 | ug/kg | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.1 | 0.18 | ug/kg | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.1 | 0.17 | ug/kg | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.1 | 0.15 | ug/kg | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.1 | 0.15 | ug/kg | | |
| 79-01-6 | Trichloroethene | 0.39 | 1.0 | 0.15 | ug/kg | J | |
| 75-69-4 | Trichlorofluoromethane | ND | 5.1 | 0.26 | ug/kg | | |
| 75-01-4 | Vinyl chloride | ND | 2.1 | 0.20 | ug/kg | | |
| | m,p-Xylene | ND | 1.0 | 0.36 | ug/kg | | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.28 | ug/kg | | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.28 | ug/kg | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 1868-53-7 | Dibromofluoromethane | 99% | | 70-1 | 22% | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 96% | | 68-1 | 24% | | |
| 2037-26-5 | Toluene-D8 | 92% | | 77-1 | 25% | | |
| 1 (0, 00, 1 | | 0.001 | | 70.1 | 2004 | | |

92%

4-Bromofluorobenzene

460-00-4

J = Indicates an estimated value

72-130%

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2

59 of 1690 ACCUTEST JC7512

| | Report of Analysis | | | | | | | |
|---|---------------------------------|-----------------------|-------------------------------|-----------------|---------------------------|-----------------------|-----------------------------|--|
| Client San Lab Sam Matrix: Method: Project: | ole ID: JC751 SO - S SW84 | 2-6 oil 6 8270D | SW846 3546 Sey Avenue, Bro | ooklyn, N | Da Pe | L |)/29/15)/31/15 5.5 | |
| Run #1 Run #2 | File ID 3E77588.D | DF 1 | Analyzed 11/03/15 | By AN | Prep Date 11/03/15 | Prep Batch OP88613 | Analytical Batch E3E3376 | |
| Run #1 Run #2 | Initial Weight 30.2 g | Final V 1.0 ml | olume | | | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|----------------------------|--------|-----|-----|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 69 | 26 | ug/kg | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 170 | 31 | ug/kg | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 170 | 28 | ug/kg | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 170 | 63 | ug/kg | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 170 | 150 | ug/kg | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 170 | 66 | ug/kg | |
| 95-48-7 | 2-Methylphenol | ND | 69 | 50 | ug/kg | |
| | 3&4-Methylphenol | ND | 69 | 33 | ug/kg | |
| 88-75-5 | 2-Nitrophenol | ND | 170 | 32 | ug/kg | |
| 100-02-7 | 4-Nitrophenol | ND | 350 | 59 | ug/kg | |
| 87-86-5 | Pentachlorophenol | ND | 170 | 85 | ug/kg | |
| 108-95-2 | Phenol | ND | 69 | 26 | ug/kg | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 170 | 33 | ug/kg | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 170 | 31 | ug/kg | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 170 | 28 | ug/kg | |
| 83-32-9 | Acenaphthene | ND | 35 | 33 | ug/kg | |
| 208-96-8 | Acenaphthylene | ND | 35 | 3.6 | ug/kg | |
| 98-86-2 | Acetophenone | ND | 170 | 5.9 | ug/kg | |
| 120-12-7 | Anthracene | ND | 35 | 3.0 | ug/kg | |
| 1912-24-9 | Atrazine | ND | 69 | 14 | ug/kg | |
| 56-55-3 | Benzo(a)anthracene | ND | 35 | 6.7 | ug/kg | |
| 50-32-8 | Benzo(a)pyrene | ND | 35 | 7.4 | ug/kg | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 35 | 7.1 | ug/kg | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 35 | 10 | ug/kg | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 35 | 7.7 | ug/kg | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 69 | 7.9 | ug/kg | |
| 85-68-7 | Butyl benzyl phthalate | ND | 69 | 19 | ug/kg | |
| 92-52-4 | 1,1'-Biphenyl | ND | 69 | 6.4 | ug/kg | |
| 100-52-7 | Benzaldehyde | ND | 170 | 8.7 | ug/kg | |
| 91-58-7 | 2-Chloronaphthalene | ND | 69 | 5.0 | ug/kg | |
| 106-47-8 | 4-Chloroaniline | ND | 170 | 9.2 | ug/kg | |
| 86-74-8 | Carbazole | ND | 69 | 3.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Client Sample ID: | SB-3-5 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC7512-6 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 95.5 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|--------|------|-------|---|
| 105-60-2 | Caprolactam | ND | 69 | 22 | ug/kg | |
| 218-01-9 | Chrysene | ND | 35 | 5.6 | ug/kg | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 69 | 7.9 | ug/kg | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 69 | 14 | ug/kg | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 69 | 7.9 | ug/kg | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 69 | 6.5 | ug/kg | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 35 | 6.5 | ug/kg | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 35 | 8.9 | ug/kg | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 69 | 23 | ug/kg | |
| 123-91-1 | 1,4-Dioxane | ND | 35 | 23 | ug/kg | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 35 | 12 | ug/kg | |
| 132-64-9 | Dibenzofuran | ND | 69 | 4.8 | ug/kg | |
| 84-74-2 | Di-n-butyl phthalate | ND | 69 | 4.1 | ug/kg | |
| 117-84-0 | Di-n-octyl phthalate | ND | 69 | 4.7 | ug/kg | |
| 84-66-2 | Diethyl phthalate | ND | 69 | 4.4 | ug/kg | |
| 131-11-3 | Dimethyl phthalate | ND | 69 | 5.0 | ug/kg | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 69 | 12 | ug/kg | |
| 206-44-0 | Fluoranthene | ND | 35 | 4.2 | ug/kg | |
| 86-73-7 | Fluorene | ND | 35 | 4.1 | ug/kg | |
| 118-74-1 | Hexachlorobenzene | ND | 69 | 6.8 | ug/kg | |
| 87-68-3 | Hexachlorobutadiene | ND | 35 | 9.2 | ug/kg | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 350 | 55 | ug/kg | |
| 67-72-1 | Hexachloroethane | ND | 170 | 11 | ug/kg | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 35 | 18 | ug/kg | |
| 78-59-1 | Isophorone | ND | 69 | 6.5 | ug/kg | |
| 91-57-6 | 2-Methylnaphthalene | ND | 69 | 6.5 | ug/kg | |
| 88-74-4 | 2-Nitroaniline | ND | 170 | 7.9 | ug/kg | |
| 99-09-2 | 3-Nitroaniline | ND | 170 | 9.8 | ug/kg | |
| 100-01-6 | 4-Nitroaniline | ND | 170 | 12 | ug/kg | |
| 91-20-3 | Naphthalene | ND | 35 | 5.5 | ug/kg | |
| 98-95-3 | Nitrobenzene | ND | 69 | 11 | ug/kg | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 69 | 10 | ug/kg | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 170 | 18 | ug/kg | |
| 85-01-8 | Phenanthrene | ND | 35 | 3.8 | ug/kg | |
| 129-00-0 | Pyrene | ND | 35 | 4.3 | ug/kg | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 170 | 8.3 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 367-12-4 | 2-Fluorophenol | 79% | | 30-1 | 06% | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

4.6 4

Page 2 of 3



Report of Analysis

| Client Sample ID: | SB-3-5 | | |
|-------------------|--|-----------------------|----------|
| Lab Sample ID: | JC7512-6 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| Method: | SW846 8270D SW846 3546 | Percent Solids: | 95.5 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN TCL List (SOM0 2.0)

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 4165-62-2 | Phenol-d5 | 76% | | 30-106% |
| 118-79-6 | 2,4,6-Tribromophenol | 87% | | 24-140% |
| 4165-60-0 | Nitrobenzene-d5 | 100% | | 26-122% |
| 321-60-8 | 2-Fluorobiphenyl | 89% | | 36-112% |
| 1718-51-0 | Terphenyl-d14 | 86% | | 36-132% |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 3 of 3

SGS

62 of 1690 ACCUTEST JC7512

| | Report of Analysis Pa | | | | | | | |
|--|------------------------------------|----------------------|-------------------------------|------------------|---------------------------|-----------------------|-----------------------------|--|
| Client Sar Lab Samı Matrix: Method: Project: | ple ID: JC7512 SO - So SW846 | -6 bil 8151 SV | W846 3550C sey Avenue, Bro | ooklyn, NY | Da Pe | ate Received: 10 |)/29/15)/31/15 5.5 | |
| Run #1 Run #2 | File ID OA112013.D | DF 2 | Analyzed 11/05/15 | By VDT | Prep Date 11/03/15 | Prep Batch OP88618 | Analytical Batch GOA3882 | |
| Run #1 Run #2 | Initial Weight 15.7 g | Final V 5.0 ml | ⁷ olume | | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|------------------|------------------|-------------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 33 6.7 6.7 | 15 2.0 2.5 | ug/kg ug/kg ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 Limits | | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2,4-DCAA | 28% 19% | | 10-1 10-1 | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



JC7512



| | Report of Analysis P. | | | | | | | |
|---|---------------------------------|---------------------------|-------------------------------|-----------------|---------------------------|-----------------------|------------------------------------|--|
| Client San Lab Sam Matrix: Method: Project: | ole ID: JC751 SO - S SW84 | 2-6 oil 6 8081B | SW846 3546 sey Avenue, Bro | ooklyn, N | Da Pe | r r | //29/15 //31/15 5.5 | |
| Run #1 Run #2 | File ID 1G116792.D | DF 1 | Analyzed 11/05/15 | By RK | Prep Date 11/03/15 | Prep Batch OP88620 | Analytical Batch G1G3831 | |
| Run #1 Run #2 | Initial Weight 16.6 g | Final V 10.0 ml | | | | | | |

Pesticide TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------|--------|--------|------|-------|---|
| 309-00-2 | Aldrin | ND | 0.63 | 0.56 | ug/kg | |
| 319-84-6 | alpha-BHC | ND | 0.63 | 0.42 | ug/kg | |
| 319-85-7 | beta-BHC | ND | 0.63 | 0.39 | ug/kg | |
| 319-86-8 | delta-BHC | ND | 0.63 | 0.25 | ug/kg | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.63 | 0.29 | ug/kg | |
| 5103-71-9 | alpha-Chlordane | ND | 0.63 | 0.34 | ug/kg | |
| 5103-74-2 | gamma-Chlordane | ND | 0.63 | 0.48 | ug/kg | |
| 60-57-1 | Dieldrin | ND | 0.63 | 0.49 | ug/kg | |
| 72-54-8 | 4,4'-DDD | ND | 0.63 | 0.23 | ug/kg | |
| 72-55-9 | 4,4'-DDE | ND | 0.63 | 0.21 | ug/kg | |
| 50-29-3 | 4,4'-DDT | ND | 0.63 | 0.24 | ug/kg | |
| 72-20-8 | Endrin | ND | 0.63 | 0.22 | ug/kg | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.63 | 0.36 | ug/kg | |
| 7421-93-4 | Endrin aldehyde | ND | 0.63 | 0.47 | ug/kg | |
| 959-98-8 | Endosulfan-I | ND | 0.63 | 0.21 | ug/kg | |
| 33213-65-9 | Endosulfan-II | ND | 0.63 | 0.60 | ug/kg | |
| 76-44-8 | Heptachlor | ND | 0.63 | 0.52 | ug/kg | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.63 | 0.26 | ug/kg | |
| 72-43-5 | Methoxychlor | ND | 1.3 | 0.35 | ug/kg | |
| 53494-70-5 | Endrin ketone | ND | 0.63 | 0.33 | ug/kg | |
| 8001-35-2 | Toxaphene | ND | 16 | 11 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 877-09-8 | Tetrachloro-m-xylene | 76% | | 24-1 | 36% | |
| 877-09-8 | Tetrachloro-m-xylene | 75% | | 24-1 | 36% | |
| 2051-24-3 | Decachlorobiphenyl | 60% | | 10-1 | 53% | |
| 2051-24-3 | Decachlorobiphenyl | 89% | | 10-1 | 53% | |

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

- E = Indicates value exceeds calibration range
- J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





| | Report of Analysis | | | | | | | |
|---|------------------------------------|--------------------------|-------------------------------|-----------------|---------------------------|-----------------------|-----------------------------|--|
| Client Sar Lab Sam Matrix: Method: Project: | ple ID: JC7512 SO - So SW846 | 2-6 pil 5 8082A | SW846 3546 sey Avenue, Bro | ooklyn, N | Da Pe | |)/29/15)/31/15 5.5 | |
| Run #1 Run #2 | File ID XX179851.D | DF 1 | Analyzed 11/07/15 | By JR | Prep Date 11/03/15 | Prep Batch OP88619 | Analytical Batch GXX5510 | |
| Run #1 Run #2 | Initial Weight 16.6 g | Final V 10.0 m | | | | | | |

PCB List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|--|--|--|---|---|---|---|
| 12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 11096-82-5 11100-14-4 37324-23-5 | Aroclor 1016 Aroclor 1221 Aroclor 1232 Aroclor 1242 Aroclor 1248 Aroclor 1254 Aroclor 1260 Aroclor 1268 Aroclor 1262 | ND ND ND ND ND ND ND ND | 32 32 32 32 32 32 32 32 32 32 32 32 | 10 19 10 14 9.7 14 13 9.7 8.9 | ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg ug/kg | |
| CAS No. 877-09-8 877-09-8 2051-24-3 2051-24-3 | Surrogate Recoveries Tetrachloro-m-xylene Tetrachloro-m-xylene Decachlorobiphenyl Decachlorobiphenyl | Run# 1 101% 99% 91% 94% | S2 8.9 ug/kj Run# 2 Limits 20-152% 20-152% 12-157% 12-157% | | 52% 52% 57% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



SGS

| Client Sample ID: | SB-3-5 | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC7512-6 | Date Sampled: | 10/29/15 |
| Matrix: | SO - Soil | Date Received: | 10/31/15 |
| | | Percent Solids: | 95.5 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|---------|-------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | 5150 | 52 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Antimony | < 2.1 | 2.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Arsenic | < 2.1 | 2.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C 2 | SW846 3050B ³ |
| Barium | 23.8 | 21 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C 2 | SW846 3050B ³ |
| Beryllium | 0.32 | 0.21 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C 2 | SW846 3050B ³ |
| Cadmium | < 0.52 | 0.52 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C 2 | SW846 3050B ³ |
| Calcium | 1210 | 520 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Chromium | 15.7 | 1.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Cobalt | 6.7 | 5.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Copper | 9.4 | 2.6 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Iron | 10600 | 52 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Lead | 4.9 | 2.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Magnesium | 2490 | 520 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Manganese | 265 | 1.6 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Mercury | < 0.031 | 0.031 | mg/kg | 1 | 11/04/15 | 11/04/15 ма | SW846 7471B ¹ | SW846 7471B ⁴ |
| Nickel | 46.1 | 4.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Potassium | < 1000 | 1000 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Selenium | < 2.1 | 2.1 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Silver | < 0.52 | 0.52 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Sodium | < 1000 | 1000 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Thallium | < 1.0 | 1.0 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Vanadium | 16.1 | 5.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |
| Zinc | 20.3 | 5.2 | mg/kg | 1 | 11/03/15 | 11/07/15 мs | SW846 6010C ² | SW846 3050B ³ |

(1) Instrument QC Batch: MA37945

(2) Instrument QC Batch: MA37978

(3) Prep QC Batch: MP90046

(4) Prep QC Batch: MP90074

RL = Reporting Limit





JC7512



Section 4

Sample Results

Report of Analysis





| | | | 1000 | | 11141.515 | | ruge r or i |
|-----------|---------------|------------|-----------------|----------|----------------|----------------|------------------|
| Client Sa | mple ID: MW | -1I-20' | | | | | |
| Lab Sam | ple ID: JC12 | 861-1 | | | Date | e Sampled: 01 | 1/18/16 |
| Matrix: | SO - | Soil | | | Date | e Received: 01 | 1/19/16 |
| Method: | SW8 | 46 8260C | SW846 5035 | | Perc | ent Solids: 82 | 2.4 |
| Project: | 2002 | -2024 Crop | sey Avenue, Bro | oklyn, N | IΥ | | |
| | | | | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | Y160757.D | 1 | 01/24/16 | PS | 01/20/16 08:00 | n/a | VY7020 |
| Run #2 | | | | | | | |
| | Initial Weigl | nt | | | | | |
| Run #1 | 4.5 g | | | | | | |
| Run #2 | - | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 13 | 3.0 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.67 | 0.18 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.7 | 0.42 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.7 | 0.21 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.7 | 0.32 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.7 | 0.49 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 13 | 2.6 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.7 | 0.31 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.7 | 0.31 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.7 | 0.21 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.7 | 0.65 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.7 | 0.20 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.7 | 0.35 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.7 | 0.43 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.7 | 0.73 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.7 | 0.28 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.3 | 0.18 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.3 | 0.16 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.3 | 0.21 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.3 | 0.30 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.7 | 0.49 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.3 | 0.19 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.3 | 0.18 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.3 | 0.80 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.3 | 1.1 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.3 | 0.80 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.7 | 0.32 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.7 | 0.16 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.7 | 0.24 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.3 | 0.22 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.7 | 0.60 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.7 | 1.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

 $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



Page 1 of 2



7 of 266

ACCUTEST

E = Indicates value exceeds calibration range

J = Indicates an estimated value

1,2,3-Trichlorobenzene

1,2,4-Trichlorobenzene

1,1,1-Trichloroethane

1,1,2-Trichloroethane

Trichlorofluoromethane

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Trichloroethene

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

o-Xylene

87-61-6

120-82-1

71-55-6

79-00-5

79-01-6

75-69-4

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

Report of Analysis

| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | | MW-1I-20' JC12861-1 SO - Soil SW846 8260C SW846 2002-2024 Cropsey Av | | klyn, NY | | Date | Sampled: Received: ent Solids: | 01/18/16 01/19/16 82.4 |
|---|---------|--|--------|----------|-------|-------|--------------------------------------|------------------------------|
| VOA TCL | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopro | pylbenzene | ND | 2.7 | 0.14 | ug/kg | | |
| 79-20-9 | Methy | l Acetate | ND | 6.7 | 1.2 | ug/kg | | |
| 108-87-2 | Methy | lcyclohexane | ND | 2.7 | 0.31 | ug/kg | | |
| 1634-04-4 | Methy | l Tert Butyl Ether | ND | 1.3 | 0.21 | ug/kg | | |
| 108-10-1 | 4-Meth | nyl-2-pentanone(MIBK) | ND | 6.7 | 0.62 | ug/kg | | |
| 75-09-2 | Methy | lene chloride | ND | 6.7 | 1.3 | ug/kg | | |
| 100-42-5 | Styren | e | ND | 2.7 | 0.24 | ug/kg | | |
| 79-34-5 | 1,1,2,2 | 2-Tetrachloroethane | ND | 2.7 | 0.24 | ug/kg | | |
| 127-18-4 | | | | 2.7 | 0.41 | ug/kg | | |
| 108-88-3 | Toluer | ne | ND | 1.3 | 0.28 | ug/kg | | |
| 0 | | | | | ~ ~ / | | | |

6.7

6.7

2.7

2.7

1.3

6.7

2.7

1.3

1.3

1.3

Run# 2

0.24

0.23

0.20

0.20

0.20

0.34

0.27

0.47

0.37

0.37

Limits

70-122%

68-124%

77-125%

72-130%

ug/kg

J

ND

ND

ND

ND

0.74

ND

ND

ND

ND

ND

Run#1

109%

110%

99%

115%

| ND = Not detected | MDL = Method Detection Limit |
|--------------------------|------------------------------|
| RL = Reporting Limit | |
| E = Indicates value exce | eeds calibration range |

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



8 of 266

ACCUTEST JC12861

Page 2 of 2

4.1 **4**



| | | | | | J >> | | ruge rorr |
|-----------|---------------|-----------|-----------------|-----------|----------------|----------------|------------------|
| Client Sa | mple ID: MW- | 1I-45' | | | | | |
| Lab Sam | ple ID: JC128 | 861-2 | | | Date | Sampled: 0 | 1/18/16 |
| Matrix: | SO - 1 | Soil | | | Date | Received: 0 | 1/19/16 |
| Method: | SW84 | 46 8260C | SW846 5035 | | Perc | ent Solids: 82 | 2.4 |
| Project: | 2002- | 2024 Crop | sey Avenue, Bro | ooklyn, N | Y | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | Y160758.D | 1 | 01/24/16 | PS | 01/20/16 08:00 | n/a | VY7020 |
| Run #2 | | | | | | | |
| | Initial Weigh | t | | | | | |
| Run #1 | 5.1 g | | | | | | |
| Run #2 | - | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 12 | 2.7 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.59 | 0.16 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.9 | 0.37 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.4 | 0.19 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.9 | 0.28 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.9 | 0.43 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 12 | 2.3 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.4 | 0.27 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.4 | 0.27 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.4 | 0.18 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.9 | 0.57 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.4 | 0.18 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.9 | 0.31 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.4 | 0.38 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.4 | 0.65 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.4 | 0.24 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.2 | 0.16 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.2 | 0.15 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.2 | 0.19 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.2 | 0.27 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.9 | 0.43 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.2 | 0.17 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.2 | 0.16 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.2 | 0.70 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.2 | 0.93 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.2 | 0.71 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.4 | 0.28 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.4 | 0.14 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.4 | 0.21 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.2 | 0.19 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.9 | 0.53 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.9 | 1.6 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



9 of 266

JC12861

Page 1 of 2

4.2

E = Indicates value exceeds calibration range

J = Indicates an estimated value

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

o-Xylene

| Report of | ² Analysis |
|------------------|-----------------------|
|------------------|-----------------------|

| Client Samp Lab Sample Matrix: Method: Project: | e ID: JC12861-2 SO - Soil SW846 8260C SW846 | JC12861-2 | | | | | 01/18/16 01/19/16 82.4 |
|---|---|-----------|------|-------|-------|---|------------------------------|
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 2.4 | 0.13 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 5.9 | 1.0 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 2.4 | 0.27 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.2 | 0.18 | ug/kg | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.9 | 0.55 | ug/kg | | |
| 75-09-2 | Methylene chloride | ND | 5.9 | 1.2 | ug/kg | | |
| 100-42-5 | Styrene | ND | 2.4 | 0.21 | ug/kg | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.4 | 0.21 | ug/kg | | |
| 127-18-4 | Tetrachloroethene | 4.3 | 2.4 | 0.36 | ug/kg | | |
| 108-88-3 | Toluene | ND | 1.2 | 0.25 | ug/kg | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.9 | 0.21 | ug/kg | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.9 | 0.20 | ug/kg | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.4 | 0.18 | ug/kg | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.4 | 0.17 | ug/kg | | |
| 79-01-6 | Trichloroethene | 0.26 | 1.2 | 0.17 | ug/kg | J | |
| 75-69-4 | ND | 5.9 | 0.30 | ug/kg | | | |

2.4

1.2

1.2

1.2

Run# 2

0.23

0.42

0.33

0.33

Limits

70-122%

68-124%

77-125%

72-130%

ug/kg

ug/kg

ug/kg

ug/kg

ND

ND

ND

ND

Run#1

107%

105%

98%

115%

| ND = Not detected | MDL = Method Detection Limit |
|--------------------------|------------------------------|
| RL = Reporting Limit | |
| E = Indicates value exce | eeds calibration range |

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



JC12861

SGS



| | Report of Analysis Pag | | | | | | | | | |
|------------------------|-------------------------------|-----------------------|------------------|------------|----------------|-------------------|------------------|--|--|--|
| Client Sar Lab Samp | - | AW-1I-50' C12861-3 | | | Date | Sampled: (|)1/18/16 | | | |
| Matrix: | S | SO - Soil | | | Date | Received: (| 01/19/16 | | | |
| Method: | S | SW846 8260C | SW846 5035 | | Perc | ent Solids: 8 | 32.4 | | | |
| Project: | 2 | 2002-2024 Crop | osey Avenue, Bro | ooklyn, NY | | | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | | | |
| Run #1 | Y160759 | .D 1 | 01/24/16 | PS | 01/20/16 08:00 | n/a | VY7020 | | | |
| Run #2 | | | | | | | | | | |
| | Initial W | eight | | | | | | | | |
| Run #1 | 3.8 g | - | | | | | | | | |
| Run #2 | - | | | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 5.7 | 16 | 3.6 | ug/kg | J |
| 71-43-2 | Benzene | ND | 0.80 | 0.21 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 8.0 | 0.49 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 3.2 | 0.25 | ug/kg | |
| 75-25-2 | Bromoform | ND | 8.0 | 0.38 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 8.0 | 0.58 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 16 | 3.0 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 3.2 | 0.36 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 3.2 | 0.37 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 3.2 | 0.25 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 8.0 | 0.77 | ug/kg | |
| 67-66-3 | Chloroform | ND | 3.2 | 0.24 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 8.0 | 0.42 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 3.2 | 0.50 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 3.2 | 0.87 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 3.2 | 0.33 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.6 | 0.21 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.6 | 0.19 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.6 | 0.25 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.6 | 0.36 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 8.0 | 0.58 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.6 | 0.23 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.6 | 0.21 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.6 | 0.95 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.6 | 1.2 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.6 | 0.95 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 3.2 | 0.38 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 3.2 | 0.19 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 3.2 | 0.28 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.6 | 0.26 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 8.0 | 0.72 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 8.0 | 2.1 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



11 of 266

JC12861

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Trichlorofluoromethane

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

o-Xylene

75-69-4

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

17060-07-0

2037-26-5

460-00-4

| Report of Analysis | S | |
|---------------------------|---|--|
|---------------------------|---|--|

| Client Samj Lab Sample Matrix: Method: Project: | - | | klyn, NY | | Date | Sampled: Received: ent Solids: | 01/18/16 01/19/16 82.4 |
|---|----------------------------|--------|----------|------|-------|--------------------------------------|------------------------------|
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 3.2 | 0.17 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 8.0 | 1.4 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 3.2 | 0.36 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.6 | 0.24 | ug/kg | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 8.0 | 0.73 | ug/kg | | |
| 75-09-2 | Methylene chloride | ND | 8.0 | 1.6 | ug/kg | | |
| 100-42-5 | Styrene | ND | 3.2 | 0.28 | ug/kg | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 3.2 | 0.28 | ug/kg | | |
| 127-18-4 | Tetrachloroethene | ND | 3.2 | 0.48 | ug/kg | | |
| 108-88-3 | Toluene | ND | 1.6 | 0.33 | ug/kg | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 8.0 | 0.28 | ug/kg | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 8.0 | 0.27 | ug/kg | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 3.2 | 0.24 | ug/kg | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 3.2 | 0.23 | ug/kg | | |
| 79-01-6 | Trichloroethene | ND | 1.6 | 0.23 | ug/kg | | |

8.0

3.2

1.6

1.6

1.6

Run# 2

0.40

0.31

0.56

0.44

0.44

Limits

70-122%

68-124%

77-125%

72-130%

ug/kg

ug/kg

ug/kg

ug/kg

ug/kg

ND

ND

ND

ND

ND

Run#1

106%

106%

98%

116%

ND = Not detectedMDL = Method Detection Limit RL = Reporting Limit E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 2



12 of 266

| Client Sample ID: | | | | | | | |
|-------------------|-------------------|-------------|-----------|----|----------------|------|--------|
| Lab Sample ID: | JC12861-4 | | | | Date Sampled: | : 01 | /18/16 |
| Matrix: | SO - Soil | | | | Date Received | : 01 | /19/16 |
| | | | | | Percent Solids | : 82 | .4 |
| Project: | 2002-2024 Cropsey | Avenue, Bro | oklyn, NY | | | | |
| | | | | | | | |
| General Chemistry | | | | | | | |
| Analyte | Result | RL | Units | DF | Analyzed | By | Method |

Report of Analysis

Page 1 of 1

4.4



| | | | Repo | ort of . | Analysis | | Page 1 of 2 |
|---|----------------------------|---|----------------------------|-----------------|-------------------------|--|------------------------------------|
| Client Sar Lab Sam Matrix: Method: Project: | ple ID: J A S | RIP C12861-5 AQ - Trip Blan W846 8260C 002-2024 Cro | k Soil psey Avenue, Bro | ooklyn, N | I I | Date Sampled: 0 Date Received: 0 Percent Solids: n | |
| Run #1 Run #2 | File ID 2A165017 | DF 7.D 1 | Analyzed 01/22/16 | By TK | Prep Date n/a | Prep Batch n/a | Analytical Batch V2A7019 |
| Run #1 Run #2 | Purge Vo 5.0 ml | lume | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.3 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.24 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.37 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.23 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.23 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.42 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 5.6 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.25 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.22 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.34 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.19 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.41 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.28 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.99 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.23 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.27 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.90 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.17 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.18 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.51 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.27 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.65 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.39 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.21 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.27 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 0.52 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.7 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



14 of 266

JC12861

E = Indicates value exceeds calibration range

Report of Analysis

| Client Sample ID: | TRIP | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC12861-5 | Date Sampled: | 01/19/16 |
| Matrix: | AQ - Trip Blank Soil | Date Received: | 01/19/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.23 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.9 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.22 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.24 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.0 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.73 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.21 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.40 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.16 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.25 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.22 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.43 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.15 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.38 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.17 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.17 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 102% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 105% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 99% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 102% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



15 of 266

JC12861

Page 2 of 2

4.5 4

| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|--|------------------------------|---------------------------------------|---------------------------|-----------------|-------------------------|--|------------------------------------|
| Client San Lab Samj Matrix: Method: Project: | ple ID: JC12 AQ - SW84 | - 861-6 Field Blanl 46 8260C | k Soil sey Avenue, Bro | ooklyn, N |] | Date Sampled: (Date Received: (Percent Solids: 1 | |
| Run #1 Run #2 | File ID 2A165018.D | DF 1 | Analyzed 01/22/16 | By TK | Prep Date n/a | Prep Batch n/a | Analytical Batch V2A7019 |
| Run #1 Run #2 | Purge Volum 5.0 ml | e | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.3 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.24 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.37 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.23 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.23 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.42 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 5.6 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.25 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.22 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.34 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.19 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.41 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.28 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.99 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.23 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.27 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.90 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.17 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.18 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.51 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.27 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.65 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.39 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.21 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.27 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 0.52 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.7 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



16 of 266

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Client Sample ID: | FIELD | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC12861-6 | Date Sampled: | 01/19/16 |
| Matrix: | AQ - Field Blank Soil | Date Received: | 01/19/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.23 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.9 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.22 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.24 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.0 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.73 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.21 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.40 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.16 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.25 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.22 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.43 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.15 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.38 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.17 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.17 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 100% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 105% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 100% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 101% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

SGS

17 of 266

ACCUTEST JC12861

4.6 **4**

Page 2 of 2

| | | | - L | | J | | | |
|-----------|--------------|------------|-----------------|-----------|----------------|----------------------|------------------|--|
| Client Sa | mple ID: MW | -45-23.5' | | | | | | |
| Lab Sam | ple ID: JC12 | 2861-7 | | | Date | Sampled: 01 | 1/19/16 | |
| Matrix: | SO - | Soil | | | Date | Received: 01 | 1/19/16 | |
| Method: | SW8 | 46 8260C | SW846 5035 | | Perc | Percent Solids: 78.7 | | |
| Project: | 2002 | -2024 Crop | sey Avenue, Bro | ooklyn, N | ΙY | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| Run #1 | Y160764.D | 1 | 01/24/16 | PS | 01/20/16 08:00 | n/a | VY7020 | |
| Run #2 | | | | | | | | |
| | Initial Weig | nt | | | | | | |
| Run #1 | 4.7 g | | | | | | | |
| Run #2 | - | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 23.0 | 14 | 3.0 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.68 | 0.18 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.8 | 0.42 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.7 | 0.21 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.8 | 0.32 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.8 | 0.49 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 14 | 2.6 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.7 | 0.31 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.7 | 0.31 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.7 | 0.21 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.8 | 0.65 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.7 | 0.20 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.8 | 0.35 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.7 | 0.43 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.7 | 0.74 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.7 | 0.28 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.4 | 0.18 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.4 | 0.16 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.4 | 0.21 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.4 | 0.30 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.8 | 0.49 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.4 | 0.19 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.4 | 0.18 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.4 | 0.80 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.4 | 1.1 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.4 | 0.80 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.7 | 0.32 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.7 | 0.16 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.7 | 0.24 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.4 | 0.22 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.8 | 0.61 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.8 | 1.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

 $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



Page 1 of 2

SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Lab Sample ID: | JC12861-7 | Date Sampled: | 01/19/16 |
|----------------|--|-----------------|----------|
| Matrix: | SO - Soil | Date Received: | 01/19/16 |
| Method: | SW846 8260C SW846 5035 | Percent Solids: | 78.7 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 2.7 | 0.14 | ug/kg | |
| 79-20-9 | Methyl Acetate | ND | 6.8 | 1.2 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 2.7 | 0.31 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.4 | 0.21 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.8 | 0.62 | ug/kg | |
| 75-09-2 | Methylene chloride | ND | 6.8 | 1.3 | ug/kg | |
| 100-42-5 | Styrene | ND | 2.7 | 0.24 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.7 | 0.24 | ug/kg | |
| 127-18-4 | Tetrachloroethene | ND | 2.7 | 0.41 | ug/kg | |
| 108-88-3 | Toluene | ND | 1.4 | 0.28 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.8 | 0.24 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.8 | 0.23 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.7 | 0.20 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.7 | 0.20 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 1.4 | 0.20 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 6.8 | 0.34 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 2.7 | 0.27 | ug/kg | |
| | m,p-Xylene | ND | 1.4 | 0.48 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 1.4 | 0.37 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 1.4 | 0.37 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 108% | | 70-1 | 22% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 105% | | 68-1 | 24% | |
| 2037-26-5 | Toluene-D8 | 99% | | 77-1 | 25% | |
| 460-00-4 | 4-Bromofluorobenzene | 116% | | 72-1 | 30% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2

SGS

| | | | | | | | 1 460 1 01 2 |
|-----------|---------------|-------------|-----------------|------------|----------------|---------------|------------------|
| Client Sa | mple ID: MW | -45-30' | | | | | |
| Lab Sam | ple ID: JC12 | 2861-8 | | | Date | Sampled: 0 | 1/19/16 |
| Matrix: | SO - | Soil | | | Date | Received: 0 | 1/19/16 |
| Method: | SW8 | 46 8260C | SW846 5035 | | Perc | ent Solids: 7 | 8.7 |
| Project: | 2002 | 2-2024 Crop | sey Avenue, Bro | ooklyn, NY | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | Y160760.D | 1 | 01/24/16 | PS | 01/20/16 08:00 | n/a | VY7020 |
| Run #2 | | | | | | | |
| | Initial Weigh | nt | | | | | |
| Run #1 | 3.9 g | | | | | | |
| Run #2 | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 58.1 | 16 | 3.6 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.81 | 0.22 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 8.1 | 0.50 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 3.3 | 0.25 | ug/kg | |
| 75-25-2 | Bromoform | ND | 8.1 | 0.38 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 8.1 | 0.59 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 16 | 3.1 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 3.3 | 0.37 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 3.3 | 0.37 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 3.3 | 0.25 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 8.1 | 0.78 | ug/kg | |
| 67-66-3 | Chloroform | ND | 3.3 | 0.24 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 8.1 | 0.43 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 3.3 | 0.51 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 3.3 | 0.89 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 3.3 | 0.33 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.6 | 0.21 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.6 | 0.20 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.6 | 0.26 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.6 | 0.37 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 8.1 | 0.59 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.6 | 0.23 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.6 | 0.22 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.6 | 0.96 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.6 | 1.3 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.6 | 0.97 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 3.3 | 0.39 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 3.3 | 0.19 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 3.3 | 0.29 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.6 | 0.27 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 8.1 | 0.73 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 8.1 | 2.2 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

20 of 266

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| | Method: SW846 8260C SW846 5035 Percent Solids: 78.7 Project: 2002-2024 Cropsey Avenue, Brooklyn, NY Percent Solids: 78.7 VOA TCL List CAS No. Compound Result RL MDL Units Q 98-82-8 Isopropylbenzene ND 3.3 0.17 ug/kg | | | | | | | |
|--------------------|---|----------|------------|-------------|----------------|---|--|--|
| VOA TCL | List | | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | | |
| 98-82-8 79-20-9 | Isopropylbenzene Methyl Acetate | ND ND | 3.3 8.1 | 0.17 1.4 | ug/kg ug/kg | | | |

| 98-82-8 | Isopropylbenzene | ND | 3.3 | 0.17 | ug/kg |
|------------|----------------------------|--------|--------|-------|-------|
| 79-20-9 | Methyl Acetate | ND | 8.1 | 1.4 | ug/kg |
| 108-87-2 | Methylcyclohexane | ND | 3.3 | 0.37 | ug/kg |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.6 | 0.25 | ug/kg |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 8.1 | 0.75 | ug/kg |
| 75-09-2 | Methylene chloride | ND | 8.1 | 1.6 | ug/kg |
| 100-42-5 | Styrene | ND | 3.3 | 0.29 | ug/kg |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 3.3 | 0.29 | ug/kg |
| 127-18-4 | Tetrachloroethene | ND | 3.3 | 0.49 | ug/kg |
| 108-88-3 | Toluene | ND | 1.6 | 0.34 | ug/kg |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 8.1 | 0.29 | ug/kg |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 8.1 | 0.28 | ug/kg |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 3.3 | 0.24 | ug/kg |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 3.3 | 0.24 | ug/kg |
| 79-01-6 | Trichloroethene | ND | 1.6 | 0.24 | ug/kg |
| 75-69-4 | Trichlorofluoromethane | ND | 8.1 | 0.41 | ug/kg |
| 75-01-4 | Vinyl chloride | ND | 3.3 | 0.32 | ug/kg |
| | m,p-Xylene | ND | 1.6 | 0.57 | ug/kg |
| 95-47-6 | o-Xylene | ND | 1.6 | 0.45 | ug/kg |
| 1330-20-7 | Xylene (total) | ND | 1.6 | 0.45 | ug/kg |
| CAS No. | Surragata Dagayanias | Run# 1 | Run# 2 | Limi | ta |
| CAS NO. | Surrogate Recoveries | KUN# 1 | KUN# 2 | LIMI | lS |
| 1868-53-7 | Dibromofluoromethane | 106% | | 70-12 | 22% |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 104% | | 68-12 | 24% |
| 2037-26-5 | Toluene-D8 | 98% | | 77-12 | 25% |
| 460-00-4 | 4-Bromofluorobenzene | 116% | | 72-13 | 30% |
| | | | | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4.8 **4**

Page 2 of 2

21 of 266

| Report of Analysis Page 1 | | | | | | | | |
|---------------------------|-----------|---------------|------------------|------------|----------------|-------------------|------------------|--|
| Client Sa | - | MW-45-40' | | | | | | |
| Lab Sam | ple ID: J | JC12861-9 | | | Date | Sampled: 0 | 1/19/16 | |
| Matrix: | | SO - Soil | | | Date | Received: 0 | 1/19/16 | |
| Method: | | SW846 8260C | SW846 5035 | | Perc | ent Solids: 7 | 8.7 | |
| Project: | | 2002-2024 Cro | psey Avenue, Bro | ooklyn, NY | Ŷ | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| Run #1 | Y160765 | .D 1 | 01/24/16 | PS | 01/20/16 08:00 | n/a | VY7020 | |
| Run #2 | | | | | | | | |
| | Initial W | eight | | | | | | |
| Run #1 | 6.7 g | | | | | | | |
| Run #2 | - | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 5.7 | 9.5 | 2.1 | ug/kg | J |
| 71-43-2 | Benzene | ND | 0.47 | 0.13 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 4.7 | 0.29 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 1.9 | 0.15 | ug/kg | |
| 75-25-2 | Bromoform | ND | 4.7 | 0.22 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 4.7 | 0.35 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 9.5 | 1.8 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 1.9 | 0.22 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 1.9 | 0.22 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 1.9 | 0.15 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 4.7 | 0.46 | ug/kg | |
| 67-66-3 | Chloroform | ND | 1.9 | 0.14 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 4.7 | 0.25 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 1.9 | 0.30 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 1.9 | 0.52 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 1.9 | 0.19 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 0.95 | 0.12 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 0.95 | 0.12 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 0.95 | 0.15 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 0.95 | 0.21 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 4.7 | 0.34 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 0.95 | 0.13 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 0.95 | 0.13 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 0.95 | 0.56 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 0.95 | 0.74 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 0.95 | 0.56 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.9 | 0.23 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.9 | 0.11 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.9 | 0.17 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 0.95 | 0.15 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 4.7 | 0.42 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 4.7 | 1.3 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





22 of 266

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Report | of | Analysis |
|--------|----|----------|
| LUPULU | | |

| | Method: SW846 8260C SW84 | | | klyn, NY | | Date | Sampled: Received: ent Solids: | 01/19/16 01/19/16 78.7 |
|------------|--------------------------|-----------------------|--------|----------|------|-------|--------------------------------------|------------------------------|
| VOA TCL I | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopro | pylbenzene | ND | 1.9 | 0.10 | ug/kg | | |
| 79-20-9 | Methy | 1 Acetate | ND | 4.7 | 0.82 | ug/kg | | |
| 108-87-2 | Methy | lcyclohexane | ND | 1.9 | 0.22 | ug/kg | | |
| 1634-04-4 | Methy | l Tert Butyl Ether | ND | 0.95 | 0.15 | ug/kg | | |
| 108-10-1 | 4-Met | hyl-2-pentanone(MIBK) | ND | 4.7 | 0.44 | ug/kg | | |
| 75-09-2 | Methylene chloride | | ND | 4.7 | 0.93 | ug/kg | | |
| 100-42-5 | Styrene | | ND | 1.9 | 0.17 | ug/kg | | |
| 79-34-5 | 1,1,2, | 2-Tetrachloroethane | ND | 1.9 | 0.17 | ug/kg | | |
| 127-18-4 | Tetrac | hloroethene | ND | 1.9 | 0.29 | ug/kg | | |
| 108-88-3 | Toluer | ne | ND | 0.95 | 0.20 | ug/kg | | |
| 87-61-6 | 1,2,3- | Trichlorobenzene | ND | 4.7 | 0.17 | ug/kg | | |
| 120-82-1 | 1,2,4- | Trichlorobenzene | ND | 4.7 | 0.16 | ug/kg | | |
| 71-55-6 | 1,1,1- | Trichloroethane | ND | 1.9 | 0.14 | ug/kg | | |
| 79-00-5 | 1,1,2- | Trichloroethane | ND | 1.9 | 0.14 | ug/kg | | |
| 79-01-6 | Trichl | oroethene | ND | 0.95 | 0.14 | ug/kg | | |
| 75-69-4 | Trichl | orofluoromethane | ND | 4.7 | 0.24 | ug/kg | | |
| 75-01-4 | Vinyl | chloride | ND | 1.9 | 0.19 | ug/kg | | |
| | m,p-X | ylene | ND | 0.95 | 0.33 | ug/kg | | |
| 95-47-6 | o-Xyle | ene | ND | 0.95 | 0.26 | ug/kg | | |
| 1330-20-7 | Xylen | e (total) | ND | 0.95 | 0.26 | ug/kg | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | Lim | iits | | |
| 1868-53-7 | Dibro | nofluoromethane | 108% | | 70-1 | 22% | | |
| 17060-07-0 | 1,2-Di | chloroethane-D4 | 107% | | 68-1 | 24% | | |
| 2037-26-5 | Toluer | ne-D8 | 99% | | 77-1 | 25% | | |
| 1 60 00 1 | | | | | | | | |

114%

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit E = Indicates value exceeds calibration range

4-Bromofluorobenzene

460-00-4

J = Indicates an estimated value

72-130%

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



23 of 266

JC12861

Page 2 of 2

4.9

| Client Sample ID: Lab Sample ID: Matrix: Project: | JC12861- SO - Soil | 10ISTURE 10 4 Cropsey A | venue, Bro | oklyn, NY | | Date Sampled: 01/19/16 Date Received: 01/19/16 Percent Solids: 78.7 | | | |
|--|-----------------------|-------------------------------|------------|-----------|----|---|----|-------------|--|
| General Chemistry | , | | | • · | | | | | |
| Analyte | | Result | RL | Units | DF | Analyzed | By | Method | |
| Moisture, Percent | | 21.3 | | % | 1 | 01/20/16 16:00 | KP | SM2540 G-97 | |

Report of Analysis

Page 1 of 1

4.10 4





Section 4

Sample Results

Report of Analysis





| Report of Analysis P | | | | | | | | |
|--|-----------------------------|---|-------------------------------|------------|---------------------------------|--------------------------------------|------------------------------|--|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC SC | W-7-23 215742-1 D - Soil W846 8260C 002-2024 Crop | SW846 5035 sey Avenue, Bro | ooklyn, NY | Date | Sampled: Received: ent Solids: | | |
| Run #1 Run #2 | File ID I211933.D | DF 1 | Analyzed 03/10/16 | By SJM | Prep Date 03/10/16 08:00 | Prep Batc n/a | h Analytical Batch VI8494 | |
| Run #1 Run #2 | Initial We 6.2 g | ight | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units Q |
|------------|-----------------------------|--------|-----|-----|---------|
| 67-64-1 | Acetone | ND | 130 | 29 | ug/kg |
| 71-43-2 | Benzene | ND | 6.4 | 1.7 | ug/kg |
| 74-97-5 | Bromochloromethane | ND | 64 | 4.0 | ug/kg |
| 75-27-4 | Bromodichloromethane | ND | 26 | 2.0 | ug/kg |
| 75-25-2 | Bromoform | ND | 64 | 3.0 | ug/kg |
| 74-83-9 | Bromomethane | ND | 64 | 4.7 | ug/kg |
| 78-93-3 | 2-Butanone (MEK) | ND | 130 | 24 | ug/kg |
| 75-15-0 | Carbon disulfide | ND | 26 | 2.9 | ug/kg |
| 56-23-5 | Carbon tetrachloride | ND | 26 | 2.9 | ug/kg |
| 108-90-7 | Chlorobenzene | ND | 26 | 2.0 | ug/kg |
| 75-00-3 | Chloroethane | ND | 64 | 6.2 | ug/kg |
| 67-66-3 | Chloroform | ND | 26 | 1.9 | ug/kg |
| 74-87-3 | Chloromethane | ND | 64 | 3.4 | ug/kg |
| 110-82-7 | Cyclohexane | ND | 26 | 4.0 | ug/kg |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 26 | 7.0 | ug/kg |
| 124-48-1 | Dibromochloromethane | ND | 26 | 2.6 | ug/kg |
| 106-93-4 | 1,2-Dibromoethane | ND | 13 | 1.7 | ug/kg |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 13 | 1.6 | ug/kg |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 13 | 2.0 | ug/kg |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 13 | 2.9 | ug/kg |
| 75-71-8 | Dichlorodifluoromethane | ND | 64 | 4.6 | ug/kg |
| 75-34-3 | 1,1-Dichloroethane | ND | 13 | 1.8 | ug/kg |
| 107-06-2 | 1,2-Dichloroethane | ND | 13 | 1.7 | ug/kg |
| 75-35-4 | 1,1-Dichloroethene | ND | 13 | 7.6 | ug/kg |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 13 | 10 | ug/kg |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 13 | 7.6 | ug/kg |
| 78-87-5 | 1,2-Dichloropropane | ND | 26 | 3.0 | ug/kg |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 26 | 1.5 | ug/kg |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 26 | 2.3 | ug/kg |
| 100-41-4 | Ethylbenzene | ND | 13 | 2.1 | ug/kg |
| 76-13-1 | Freon 113 | ND | 64 | 5.7 | ug/kg |
| 591-78-6 | 2-Hexanone | ND | 64 | 17 | ug/kg |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



8 of 257

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Report | of | Analysis |
|--------|----|----------|
|--------|----|----------|

| Client Sample ID: | MW-7-23 | | |
|--------------------------|--|------------------------|----------|
| Lab Sample ID: | JC15742-1 | Date Sampled: | 03/08/16 |
| Matrix: | SO - Soil | Date Received: | 03/09/16 |
| Method: | SW846 8260C SW846 5035 | Percent Solids: | 6.3 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. Compound | | Result | RL | MDL | Units | Q |
|------------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 26 | 1.4 | ug/kg | |
| 79-20-9 | Methyl Acetate | ND | 64 | 11 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 26 | 2.9 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 13 | 2.0 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 64 | 5.9 | ug/kg | |
| 75-09-2 | Methylene chloride | ND | 64 | 13 | ug/kg | |
| 100-42-5 | Styrene | ND | 26 | 2.3 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 26 | 2.2 | ug/kg | |
| 127-18-4 | Tetrachloroethene | ND | 26 | 3.9 | ug/kg | |
| 108-88-3 | Toluene | ND | 13 | 2.7 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 64 | 2.3 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 64 | 2.2 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 26 | 1.9 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 26 | 1.9 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 13 | 1.9 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 64 | 3.2 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 26 | 2.5 | ug/kg | |
| | m,p-Xylene | ND | 13 | 4.5 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 13 | 3.5 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 13 | 3.5 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 97% | | 70-12 | 22% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 95% | | 68-12 | 24% | |
| 2037-26-5 | Toluene-D8 | 98% | | 77-12 | 25% | |
| 460-00-4 | 4-Bromofluorobenzene | 104% | | 72-1 | 30% | |

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound

Page 2 of 2

9 of 257

| | | | | | | | 1 4 90 1 01 2 |
|-----------|---------------|------------|-----------------|-----------|----------------|---------------|------------------|
| Client Sa | mple ID: MW- | 7-34 | | | | | |
| Lab Sam | ple ID: JC15 | 742-2 | | | Date | Sampled: 0 | 3/08/16 |
| Matrix: | SO - | Soil | | | Date | Received: 0 | 3/09/16 |
| Method: | SW8 | 46 8260C | SW846 5035 | | Perc | ent Solids: 6 | .3 |
| Project: | 2002 | -2024 Crop | sey Avenue, Bro | oklyn, NY | Y | | |
| | | DE | | D | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | I211934.D | 1 | 03/10/16 | SJM | 03/10/16 08:00 | n/a | VI8494 |
| Run #2 | | | | | | | |
| | Initial Weigh | ıt | | | | | |
| Run #1 | 5.4 g | | | | | | |
| Run #2 | - | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units Q |
|------------|-----------------------------|--------|-----|-----|---------|
| 67-64-1 | Acetone | ND | 150 | 33 | ug/kg |
| 71-43-2 | Benzene | ND | 7.3 | 2.0 | ug/kg |
| 74-97-5 | Bromochloromethane | ND | 73 | 4.5 | ug/kg |
| 75-27-4 | Bromodichloromethane | ND | 29 | 2.3 | ug/kg |
| 75-25-2 | Bromoform | ND | 73 | 3.5 | ug/kg |
| 74-83-9 | Bromomethane | ND | 73 | 5.3 | ug/kg |
| 78-93-3 | 2-Butanone (MEK) | ND | 150 | 28 | ug/kg |
| 75-15-0 | Carbon disulfide | ND | 29 | 3.4 | ug/kg |
| 56-23-5 | Carbon tetrachloride | ND | 29 | 3.4 | ug/kg |
| 108-90-7 | Chlorobenzene | ND | 29 | 2.3 | ug/kg |
| 75-00-3 | Chloroethane | ND | 73 | 7.1 | ug/kg |
| 67-66-3 | Chloroform | ND | 29 | 2.2 | ug/kg |
| 74-87-3 | Chloromethane | ND | 73 | 3.9 | ug/kg |
| 110-82-7 | Cyclohexane | ND | 29 | 4.6 | ug/kg |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 29 | 8.0 | ug/kg |
| 124-48-1 | Dibromochloromethane | ND | 29 | 3.0 | ug/kg |
| 106-93-4 | 1,2-Dibromoethane | ND | 15 | 1.9 | ug/kg |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 15 | 1.8 | ug/kg |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 15 | 2.3 | ug/kg |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 15 | 3.3 | ug/kg |
| 75-71-8 | Dichlorodifluoromethane | ND | 73 | 5.3 | ug/kg |
| 75-34-3 | 1,1-Dichloroethane | ND | 15 | 2.1 | ug/kg |
| 107-06-2 | 1,2-Dichloroethane | ND | 15 | 2.0 | ug/kg |
| 75-35-4 | 1,1-Dichloroethene | ND | 15 | 8.7 | ug/kg |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 15 | 11 | ug/kg |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 15 | 8.7 | ug/kg |
| 78-87-5 | 1,2-Dichloropropane | ND | 29 | 3.5 | ug/kg |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 29 | 1.7 | ug/kg |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 29 | 2.6 | ug/kg |
| 100-41-4 | Ethylbenzene | ND | 15 | 2.4 | ug/kg |
| 76-13-1 | Freon 113 | ND | 73 | 6.6 | ug/kg |
| 591-78-6 | 2-Hexanone | ND | 73 | 20 | ug/kg |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





Page 1 of 2



E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Report | of | Analysis |
|--------|----|----------|
|--------|----|----------|

| Client San | nple ID: MW-7-34 | | | | | | | |
|------------|-------------------|------------------------|-----------|-----------------|-------|---------------------------------|-----|--|
| Lab Samp | le ID: JC15742-2 | JC15742-2 SO - Soil | | | | Date Sampled: Date Received: | | |
| Matrix: | SO - Soil | | | | | | | |
| Method: | SW846 8260C S | W846 5035 | | Percent Solids: | | | 6.3 | |
| Project: | 2002-2024 Cropse | ey Avenue, Broo | oklyn, NY | r | | | | |
| VOA TCL | A LISU | | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | | |
| 98-82-8 | Isopropylbenzene | ND | 29 | 1.6 | ug/kg | | | |
| 79-20-9 | Methyl Acetate | ND | 73 | 13 | ug/kg | | | |
| 108-87-2 | Methylcvclohexane | ND | 29 | 3.3 | ug/kg | | | |

| 98-82-8 | Isopropylbenzene | ND | 29 | 1.6 | ug/kg | |
|------------|----------------------------|--------|--------|-------|-------|---|
| 79-20-9 | Methyl Acetate | ND | 73 | 13 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 29 | 3.3 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 15 | 2.2 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 73 | 6.7 | ug/kg | |
| 75-09-2 | Methylene chloride | ND | 73 | 14 | ug/kg | |
| 100-42-5 | Styrene | ND | 29 | 2.6 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 29 | 2.6 | ug/kg | |
| 127-18-4 | Tetrachloroethene | 7.6 | 29 | 4.4 | ug/kg | J |
| 108-88-3 | Toluene | ND | 15 | 3.1 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 73 | 2.6 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 73 | 2.5 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 29 | 2.2 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 29 | 2.2 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 15 | 2.2 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 73 | 3.7 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 29 | 2.9 | ug/kg | |
| | m,p-Xylene | ND | 15 | 5.2 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 15 | 4.0 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 15 | 4.0 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 97% | | 70-12 | 22% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 99% | | 68-12 | | |
| 2037-26-5 | Toluene-D8 | 98% | | 77-12 | 25% | |
| 460-00-4 | 4-Bromofluorobenzene | 101% | | 72-13 | 30% | |
| | | | | | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4.2

4

Page 2 of 2

SGS

| Report of Analysis | | | | | | | | |
|--|-----------------------------|--|-------------------------------|------------|---------------------------------|--------------------------------------|-------------------------------------|--|
| Client Sa Lab Sam Matrix: Method: Project: | ple ID: J S | 4W-6-23 C15742-3 O - Soil W846 8260C 002-2024 Crop | SW846 5035 sey Avenue, Bro | ooklyn, NY | Date | Sampled: Received: ent Solids: | | |
| Run #1 Run #2 | File ID I211935.1 | DF D 1 | Analyzed 03/10/16 | By SJM | Prep Date 03/10/16 08:00 | Prep Batcl n/a | h Analytical Batch VI8494 | |
| Run #1 Run #2 | Initial W 4.8 g | eight | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|-----|-----|-------|---|
| 67-64-1 | Acetone | 203 | 170 | 38 | ug/kg | |
| 71-43-2 | Benzene | ND | 8.4 | 2.2 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 84 | 5.2 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 34 | 2.6 | ug/kg | |
| 75-25-2 | Bromoform | ND | 84 | 4.0 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 84 | 6.1 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 170 | 32 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 34 | 3.8 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 34 | 3.9 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 34 | 2.6 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 84 | 8.1 | ug/kg | |
| 67-66-3 | Chloroform | ND | 34 | 2.5 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 84 | 4.4 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 34 | 5.3 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 34 | 9.1 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 34 | 3.4 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 17 | 2.2 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 17 | 2.0 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 17 | 2.6 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 17 | 3.8 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 84 | 6.1 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 17 | 2.4 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 17 | 2.3 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 17 | 9.9 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 17 | 13 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 17 | 10 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 34 | 4.0 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 34 | 2.0 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 34 | 3.0 | ug/kg | |
| 100-41-4 | Ethylbenzene | 10 | 17 | 2.7 | ug/kg | J |
| 76-13-1 | Freon 113 | ND | 84 | 7.5 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 84 | 23 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



12 of 257

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Report | of | Analys | sis |
|--------|----|--------|-----|
|--------|----|--------|-----|

| Client Sample ID: | MW-6-23 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC15742-3 | Date Sampled: | 03/08/16 |
| Matrix: | SO - Soil | Date Received: | 03/09/16 |
| Method: | SW846 8260C SW846 5035 | Percent Solids: | 6.2 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 34 | 1.8 | ug/kg | |
| 79-20-9 | Methyl Acetate | ND | 84 | 14 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 34 | 3.8 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 17 | 2.6 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 84 | 7.7 | ug/kg | |
| 75-09-2 | Methylene chloride | ND | 84 | 16 | ug/kg | |
| 100-42-5 | Styrene | ND | 34 | 3.0 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 34 | 2.9 | ug/kg | |
| 127-18-4 | Tetrachloroethene | ND | 34 | 5.1 | ug/kg | |
| 108-88-3 | Toluene | 143 | 17 | 3.5 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 84 | 3.0 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 84 | 2.8 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 34 | 2.5 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 34 | 2.5 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 17 | 2.5 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 84 | 4.2 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 34 | 3.3 | ug/kg | |
| | m,p-Xylene | 42.3 | 17 | 5.9 | ug/kg | |
| 95-47-6 | o-Xylene | 14.1 | 17 | 4.6 | ug/kg | J |
| 1330-20-7 | Xylene (total) | 56.4 | 17 | 4.6 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 1868-53-7 | Dibromofluoromethane | 97% | | 70-1 | 22% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 98% | | 68-1 | 24% | |
| 2037-26-5 | Toluene-D8 | 98% | | 77-1 | 25% | |
| 460-00-4 | 4-Bromofluorobenzene | 103% | | 72-1 | 30% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



JC15742

Page 2 of 2

| | | | | | | | ruge r or |
|-----------|---------------|------------|-----------------|------------|----------------|---------------|------------------|
| Client Sa | mple ID: MW | -6-30 | | | | | |
| Lab Sam | ple ID: JC15 | 742-4 | | | Date | Sampled: (| 03/08/16 |
| Matrix: | SO - | Soil | | | Date | Received: (| 03/09/16 |
| Method: | SW8 | 46 8260C | SW846 5035 | | Perc | ent Solids: 6 | 5.2 |
| Project: | 2002 | -2024 Crop | sey Avenue, Bro | ooklyn, NY | ľ | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | I211936.D | 1 | 03/10/16 | SJM | 03/10/16 08:00 | n/a | VI8494 |
| Run #2 | | | | | | | |
| | Initial Weigh | nt | | | | | |
| Run #1 | 3.6 g | | | | | | |
| Run #2 | - | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|-----|-----|-------|---|
| 67-64-1 | Acetone | 283 | 220 | 50 | ug/kg | |
| 71-43-2 | Benzene | ND | 11 | 3.0 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 110 | 6.9 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 45 | 3.5 | ug/kg | |
| 75-25-2 | Bromoform | ND | 110 | 5.3 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 110 | 8.2 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 220 | 43 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 45 | 5.1 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 45 | 5.2 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 45 | 3.5 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 110 | 11 | ug/kg | |
| 67-66-3 | Chloroform | ND | 45 | 3.3 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 110 | 5.9 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 45 | 7.1 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 45 | 12 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 45 | 4.6 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 22 | 2.9 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 22 | 2.7 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 22 | 3.5 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 22 | 5.0 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 110 | 8.1 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 22 | 3.2 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 22 | 3.0 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 22 | 13 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 22 | 17 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 22 | 13 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 45 | 5.3 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 45 | 2.6 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 45 | 4.0 | ug/kg | |
| 100-41-4 | Ethylbenzene | 8.8 | 22 | 3.7 | ug/kg | J |
| 76-13-1 | Freon 113 | ND | 110 | 10 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 110 | 30 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 1 of 2

14 of 257

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Report | of | Analys | sis |
|--------|----|--------|-----|
|--------|----|--------|-----|

| Client Samj Lab Sample Matrix: Method: Project: | SO - Soil SW846 82600 | C SW846 5035 opsey Avenue, Broo | klyn, NY | 7 | Date | Sampled: Received: ent Solids: | |
|---|--------------------------|------------------------------------|----------|-----|-------|--------------------------------------|--|
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 00.0 0 .0 | . | MD | | | 1 | | |

| 98-82-8 | Isopropylbenzene | ND | 45 | 2.4 | ug/kg | |
|------------|----------------------------|--------|--------|-------|-------|---|
| 79-20-9 | Methyl Acetate | ND | 110 | 19 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 45 | 5.1 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 22 | 3.4 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 110 | 10 | ug/kg | |
| 75-09-2 | Methylene chloride | ND | 110 | 22 | ug/kg | |
| 100-42-5 | Styrene | ND | 45 | 4.0 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 45 | 3.9 | ug/kg | |
| 127-18-4 | Tetrachloroethene | ND | 45 | 6.7 | ug/kg | |
| 108-88-3 | Toluene | 124 | 22 | 4.7 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 110 | 3.9 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 110 | 3.8 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 45 | 3.3 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 45 | 3.3 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 22 | 3.3 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 110 | 5.6 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 45 | 4.4 | ug/kg | |
| | m,p-Xylene | 33.3 | 22 | 7.9 | ug/kg | |
| 95-47-6 | o-Xylene | 10.3 | 22 | 6.1 | ug/kg | J |
| 1330-20-7 | Xylene (total) | 43.6 | 22 | 6.1 | ug/kg | |
| | | | | | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limit | ts | |
| | | | | | | |
| 1868-53-7 | Dibromofluoromethane | 95% | | 70-12 | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 99% | | 68-12 | | |
| 2037-26-5 | Toluene-D8 | 98% | | 77-12 | | |
| 460-00-4 | 4-Bromofluorobenzene | 103% | | 72-13 | 80% | |
| | | | | | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4.4

Page 2 of 2



15 of 257

JC15742

| Report of Analysis Pa | | | | | | | | | |
|---|---------------------------|-----------|---------------------------------|------------------|---------------------------------|--------------------------|------------------------------|--|--|
| Client Sa Lab Samj Matrix: Method: Project: | ple ID: | | SW846 5035 opsey Avenue, Bro | ooklyn, NY | Date | Received: | 03/08/16 03/09/16 79.7 | | |
| Run #1 Run #2 | File ID I211937 | DF D 1 | Analyzed 03/10/16 | By SJM | Prep Date 03/10/16 08:00 | Prep Batch n/a | Analytical Batch VI8494 | | |
| Run #1 Run #2 | Initial V 4.8 g | Veight | | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 12.9 | 13 | 2.9 | ug/kg | J |
| 71-43-2 | Benzene | ND | 0.65 | 0.17 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.5 | 0.40 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.6 | 0.20 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.5 | 0.31 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.5 | 0.48 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 13 | 2.5 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.6 | 0.30 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.6 | 0.30 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.6 | 0.20 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.5 | 0.63 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.6 | 0.19 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.5 | 0.34 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.6 | 0.41 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.6 | 0.71 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.6 | 0.27 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.3 | 0.17 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.3 | 0.16 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.3 | 0.21 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.3 | 0.29 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.5 | 0.47 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.3 | 0.18 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.3 | 0.18 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.3 | 0.77 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.3 | 1.0 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.3 | 0.78 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.6 | 0.31 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.6 | 0.15 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.6 | 0.23 | ug/kg | |
| 100-41-4 | Ethylbenzene | 5.3 | 1.3 | 0.21 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.5 | 0.59 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.5 | 1.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



JC15742

4.5 **4**

Compound

Isopropylbenzene

CAS No.

98-82-8

Report of Analysis

| Lab Sample ID: Matrix: | JC15742-5 SO - Soil | Date Sampled: Date Received: | |
|---------------------------|--|---------------------------------|--|
| Method: | SW846 8260C SW846 5035 | Percent Solids: | |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

RL

2.6

MDL

0.14

Units

ug/kg

Q

Result

ND

| 90-02-0 | isopiopyidenzene | ND | 2.0 | 0.14 | ug/ Kg |
|------------|----------------------------|--------|--------|-------|--------|
| 79-20-9 | Methyl Acetate | ND | 6.5 | 1.1 | ug/kg |
| 108-87-2 | Methylcyclohexane | ND | 2.6 | 0.30 | ug/kg |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.3 | 0.20 | ug/kg |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.5 | 0.60 | ug/kg |
| 75-09-2 | Methylene chloride | ND | 6.5 | 1.3 | ug/kg |
| 100-42-5 | Styrene | ND | 2.6 | 0.23 | ug/kg |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.6 | 0.23 | ug/kg |
| 127-18-4 | Tetrachloroethene | 11.4 | 2.6 | 0.39 | ug/kg |
| 108-88-3 | Toluene | 117 | 1.3 | 0.27 | ug/kg |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.5 | 0.23 | ug/kg |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.5 | 0.22 | ug/kg |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.6 | 0.19 | ug/kg |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.6 | 0.19 | ug/kg |
| 79-01-6 | Trichloroethene | ND | 1.3 | 0.19 | ug/kg |
| 75-69-4 | Trichlorofluoromethane | ND | 6.5 | 0.33 | ug/kg |
| 75-01-4 | Vinyl chloride | ND | 2.6 | 0.26 | ug/kg |
| | m,p-Xylene | 20.9 | 1.3 | 0.46 | ug/kg |
| 95-47-6 | o-Xylene | 5.6 | 1.3 | 0.36 | ug/kg |
| 1330-20-7 | Xylene (total) | 26.5 | 1.3 | 0.36 | ug/kg |
| | | | | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limit | ts |
| | | | | | |
| 1868-53-7 | Dibromofluoromethane | 98% | | 70-12 | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 101% | | 68-12 | 24% |
| 2037-26-5 | Toluene-D8 | 97% | | 77-12 | |
| 460-00-4 | 4-Bromofluorobenzene | 102% | | 72-13 | 80% |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4.5

Page 2 of 2

SGS

17 of 257

| Client San Lab Samp Matrix: Method: Project: | le ID: JC1574 SO - So SW846 | W846 5035 y Avenue, Bro | 7846 5035 Avenue, Brooklyn, NY | | | Date Sampled: 03/08/16 Date Received: 03/09/16 Percent Solids: 79.7 | | | | |
|--|-----------------------------------|----------------------------|-----------------------------------|------------|---------|---|------------|------------------|--|--|
| | File ID | DF | Analyzed | By | Prep D | | Prep Batch | Analytical Batch | | |
| Run #1 | I211938.D | 1 | 03/10/16 | SJM | | 6 08:00 | n/a | VI8494 | | |
| Run #2 | D238880.D | 1 | 03/14/16 | AJ | 03/10/1 | 6 08:00 | n/a | VD9736 | | |
| | Initial Weight | Final Vo | lume N | Iethanol A | liquot | | | | | |
| Run #1 | 4.8 g | | | | • | | | | | |
| Run #2 | 4.2 g | 5.0 ml | 1 | 00 ul | | | | | | |
| VOA TCL | List | | | | | | | | | |
| CAS No. | Compound | | Result | RL | MDL | Units | Q | | | |
| 67-64-1 | Acetone | | ND | 13 | 2.9 | ug/kg | | | | |
| 71-43-2 | Benzene | | ND | 0.65 | 0.17 | ug/kg | | | | |
| 74-97-5 | Bromochlorom | ethane | ND | 6.5 | 0.40 | ug/kg | | | | |
| 75-27-4 | Bromodichloro | methane | ND | 2.6 | 0.20 | ug/kg | | | | |
| 75-25-2 | Bromoform | | ND | 6.5 | 0.31 | ug/kg | | | | |
| 74-83-9 | Bromomethane | | ND | 6.5 | 0.48 | ug/kg | | | | |
| 78-93-3 | 2-Butanone (M | EK) | ND | 13 | 2.5 | ug/kg | | | | |

Report of Analysis

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 13 | 2.9 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.65 | 0.17 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.5 | 0.40 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.6 | 0.20 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.5 | 0.31 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.5 | 0.48 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 13 | 2.5 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.6 | 0.30 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.6 | 0.30 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.6 | 0.20 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.5 | 0.63 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.6 | 0.19 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.5 | 0.34 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.6 | 0.41 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.6 | 0.71 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.6 | 0.27 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.3 | 0.17 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.3 | 0.16 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.3 | 0.21 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.3 | 0.29 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.5 | 0.47 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.3 | 0.18 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.3 | 0.18 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.3 | 0.77 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | 2440 a | 87 | 68 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.3 | 0.78 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.6 | 0.31 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.6 | 0.15 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.6 | 0.23 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.3 | 0.21 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.5 | 0.59 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.5 | 1.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



18 of 257

E = Indicates value exceeds calibration range

| Client Sam | - | | | | Data | Somplade | 03/08/16 |
|----------------------|-----------------|------------------|----------|-----------------------|----------|-------------|----------|
| Lab Sampl Matrix: | SO - Soil | | | Sampled: Received: | 03/09/16 | | |
| Method: | | | | | | ent Solids: | 79.7 |
| Project: | 2002-2024 Crops | sey Avenue, Broo | klyn, NY | , | | | |
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| | | | | | | | |

| | I I I I I I I I I I I I I I I I I I I | | | | |
|------------|---------------------------------------|-------------------|--------|-------|-------|
| 98-82-8 | Isopropylbenzene | ND | 2.6 | 0.14 | ug/kg |
| 79-20-9 | Methyl Acetate | ND | 6.5 | 1.1 | ug/kg |
| 108-87-2 | Methylcyclohexane | ND | 2.6 | 0.30 | ug/kg |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.3 | 0.20 | ug/kg |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.5 | 0.60 | ug/kg |
| 75-09-2 | Methylene chloride | ND | 6.5 | 1.3 | ug/kg |
| 100-42-5 | Styrene | ND | 2.6 | 0.23 | ug/kg |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.6 | 0.23 | ug/kg |
| 127-18-4 | Tetrachloroethene | 2740 ^a | 170 | 26 | ug/kg |
| 108-88-3 | Toluene | ND | 1.3 | 0.27 | ug/kg |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.5 | 0.23 | ug/kg |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.5 | 0.22 | ug/kg |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.6 | 0.19 | ug/kg |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.6 | 0.19 | ug/kg |
| 79-01-6 | Trichloroethene | 101 | 1.3 | 0.19 | ug/kg |
| 75-69-4 | Trichlorofluoromethane | ND | 6.5 | 0.33 | ug/kg |
| 75-01-4 | Vinyl chloride | ND | 2.6 | 0.26 | ug/kg |
| | m,p-Xylene | ND | 1.3 | 0.46 | ug/kg |
| 95-47-6 | o-Xylene | ND | 1.3 | 0.36 | ug/kg |
| 1330-20-7 | Xylene (total) | ND | 1.3 | 0.36 | ug/kg |
| | | | | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts |
| | | 0.501 | 1050 | 50.40 | |
| 1868-53-7 | Dibromofluoromethane | 95% | 107% | 70-12 | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 98% | 109% | 68-12 | |
| 2037-26-5 | Toluene-D8 | 97% | 99% | 77-12 | |
| 460-00-4 | 4-Bromofluorobenzene | 102% | 98% | 72-13 | 60% |
| | | | | | |

(a) Result is from Run# 2

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



JC15742

4.6 **4**

| Client San Lab Samp Matrix: Method: Project: | ole ID: JC1574 SO - So SW846 | 2-7 bil 8260C S | W846 5035 y Avenue, Bro | ooklyn, NY | | Date | Received: 03 | 8/08/16 8/09/16 9.7 |
|--|------------------------------------|-----------------------|----------------------------|------------|---------|----------|--------------|---------------------------|
| | File ID | DF | Analyzed | By | Prep D | ate | Prep Batch | Analytical Batch |
| Run #1 | I211939.D | 1 | 03/10/16 | SJM | 03/10/1 | 16 08:00 | n/a | VI8494 |
| Run #2 | D238881.D | 1 | 03/14/16 | AJ | 03/10/1 | 16 08:00 | n/a | VD9736 |
| | Initial Weight | Final Vo | olume 1 | Methanol A | liquot | | | |
| Run #1 | 4.8 g | | | | | | | |
| Run #2 | 4.0 g | 5.0 ml | - | 100 ul | | | | |
| VOA TCI | L List | | | | | | | |
| CAS No. | Compound | | Result | RL | MDL | Units | Q | |
| 67-64-1 | Acetone | | ND | 13 | 2.9 | ug/kg | | |
| 71-43-2 | Benzene | | ND | 0.65 | 0.17 | ug/kg | | |
| 74-97-5 | Bromochlorom | ethane | ND | 6.5 | 0.40 | ug/kg | | |
| 75-27-4 | Bromodichloro | methane | ND | 2.6 | 0.20 | ug/kg | | |

Report of Analysis

Page 1 of 2

4.7

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|-------------------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 13 | 2.9 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.65 | 0.17 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.5 | 0.40 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.6 | 0.20 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.5 | 0.31 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.5 | 0.48 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 13 | 2.5 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.6 | 0.30 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.6 | 0.30 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.6 | 0.20 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.5 | 0.63 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.6 | 0.19 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.5 | 0.34 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.6 | 0.41 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.6 | 0.71 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.6 | 0.27 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.3 | 0.17 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.31 | 1.3 | 0.16 | ug/kg | J |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.3 | 0.21 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.3 | 0.29 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.5 | 0.47 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.3 | 0.18 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.3 | 0.18 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | 2.5 | 1.3 | 0.77 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | 3670 ^a | 91 | 71 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | 4.4 | 1.3 | 0.78 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.6 | 0.31 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.6 | 0.15 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.6 | 0.23 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.3 | 0.21 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.5 | 0.59 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.5 | 1.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



| Client Sample ID: | MW-5-35 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC15742-7 | Date Sampled: | 03/08/16 |
| Matrix: | SO - Soil | Date Received: | 03/09/16 |
| Method: | SW846 8260C SW846 5035 | Percent Solids: | 79.7 |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|-------------------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | 0.15 | 2.6 | 0.14 | ug/kg | J |
| 79-20-9 | Methyl Acetate | ND | 6.5 | 1.1 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 2.6 | 0.30 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.3 | 0.20 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.5 | 0.60 | ug/kg | |
| 75-09-2 | Methylene chloride | ND | 6.5 | 1.3 | ug/kg | |
| 100-42-5 | Styrene | ND | 2.6 | 0.23 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.6 | 0.23 | ug/kg | |
| 127-18-4 | Tetrachloroethene | 3350 ^a | 180 | 27 | ug/kg | |
| 108-88-3 | Toluene | ND | 1.3 | 0.27 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.5 | 0.23 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.5 | 0.22 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.6 | 0.19 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.6 | 0.19 | ug/kg | |
| 79-01-6 | Trichloroethene | 636 ^a | 91 | 13 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 6.5 | 0.33 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 2.6 | 0.26 | ug/kg | |
| | m,p-Xylene | ND | 1.3 | 0.46 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 1.3 | 0.36 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 1.3 | 0.36 | ug/kg | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 97% | 105% | 70-12 | 22% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 97% | 109% | 68-12 | 24% | |
| 2037-26-5 | Toluene-D8 | 98% | 97% | 77-12 | 25% | |
| 460-00-4 | 4-Bromofluorobenzene | 103% | 99% | 72-13 | 30% | |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2

JC15742

| Client Sample ID: | MW-5 | | | | | | |
|------------------------------|---------------------|-------------|-----------|----|----------------|------|--------|
| Lab Sample ID: | JC15742-8 | | | | Date Sampled | : 03 | /09/16 |
| Matrix: | SO - Soil | | | | Date Received | : 03 | /09/16 |
| | | | | | Percent Solids | : 79 | .7 |
| Project: | 2002-2024 Cropsey A | Avenue, Bro | oklyn, NY | | | | |
| | | | | | | | |
| General Chemistry | 7 | | | | | | |
| General Chemistry Analyte | Result | RL | Units | DF | Analyzed | By | Method |

Report of Analysis

Page 1 of 1

4.8





Section 4

Sample Results

Report of Analysis





| Lab Sam Matrix: Method: Project: | ple ID: JC AC SV | RIP BLANK 16088-1 Q - Trip Blank V846 8260C 02-2024 Crop | soil sey Avenue, Bro | ooklyn, NY | Da | ate Sampled: 03 ate Received: 03 ercent Solids: n/ | |
|---|-----------------------------|--|--------------------------|-----------------|-------------------------|--|----------------------------|
| Run #1 Run #2 | File ID U204157.I | DF D 1 | Analyzed 03/15/16 | By NH | Prep Date n/a | Prep Batch n/a | Analytical Batch VU9383 |
| Run #1 Run #2 | Purge Vol 5.0 ml | ume | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.3 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.24 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.37 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.23 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.23 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.42 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 5.6 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.25 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.22 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.34 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.19 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.41 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.28 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.99 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.23 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.27 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.90 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.17 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.18 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.51 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.27 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.65 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.39 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.21 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.27 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 0.52 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.7 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $J = \ Indicates \ an \ estimated \ value$

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

SGS

E = Indicates value exceeds calibration range

| Client Sample ID: | TRIP BLANK | | |
|-------------------|--|-----------------------|----------|
| Lab Sample ID: | JC16088-1 | Date Sampled: | 03/10/16 |
| Matrix: | AQ - Trip Blank Soil | Date Received: | 03/11/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.23 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.9 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.22 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.24 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.0 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.73 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.21 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.40 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.16 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.25 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.22 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.43 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.15 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.38 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.17 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.17 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 105% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 107% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 100% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 104% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



SGS

8 of 128

ACCUTEST JC16088

Page 2 of 2

4.1 **4**

| Client Sa Lab Sam Matrix: Method: Project: | ple ID: JC1 AQ SW | FIELD BLANK JC16088-2 AQ - Field Blank Soil SW846 8260C 2002-2024 Cropsey Avenue, Brooklyn, NY | | | Da | ate Sampled: 03 ate Received: 03 ercent Solids: n/ | | |
|--|-----------------------------|--|--------------------------|-----------------|-------------------------|--|-----------------------------------|--|
| Run #1 Run #2 | File ID U204158.D | DF 1 | Analyzed 03/15/16 | By NH | Prep Date n/a | Prep Batch n/a | Analytical Batch VU9383 | |
| Run #1 Run #2 | Purge Volu 5.0 ml | me | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.3 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.24 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.37 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.23 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.23 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.42 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 5.6 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.25 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.22 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.34 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.19 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.41 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.28 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.99 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.23 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.27 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.90 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.17 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.18 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.51 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.27 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.65 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.39 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.21 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.27 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 0.52 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.7 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

9 of 128

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Sample ID: | FIELD BLANK | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC16088-2 | Date Sampled: | 03/10/16 |
| Matrix: | AQ - Field Blank Soil | Date Received: | 03/11/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.23 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.9 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.22 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.24 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.0 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.73 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.21 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.40 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.16 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.25 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.22 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.43 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.15 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.38 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.17 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.17 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 105% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 107% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 99% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 105% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





10 of 128

| Client Sample ID: | MW-6 | | | | | | |
|------------------------------|-------------------|--|-------|----|----------------|------|--------|
| Lab Sample ID: | JC16088-3 | | | | Date Sampled | : 03 | /10/16 |
| Matrix: | SO - Soil | | | | Date Received | : 03 | /11/16 |
| | | | | | Percent Solids | : 93 | .8 |
| Project: | 2002-2024 Cropsey | 2002-2024 Cropsey Avenue, Brooklyn, NY | | | | | |
| | | | | | | | |
| General Chemistry | | | | | | | |
| General Chemistry Analyte | Result | RL | Units | DF | Analyzed | By | Method |

Report of Analysis

Page 1 of 1

4.3



| Client Sample ID: | MW-7 | | | | | | |
|--------------------------|-------------------|-------------|-----------|----|----------------|------|--------|
| Lab Sample ID: | JC16088-4 | | | | Date Sampled | : 03 | /10/16 |
| Matrix: | SO - Soil | | | | Date Received | : 03 | /11/16 |
| | | | | | Percent Solids | : 93 | .7 |
| Project: | 2002-2024 Cropsey | Avenue, Bro | oklyn, NY | | | | |
| General Chemistry | , | | | | | | |
| | | | | | | | |
| Analyte | Result | RL | Units | DF | Analyzed | By | Method |

Report of Analysis

Page 1 of 1

4.4





Section 4

Sample Results

Report of Analysis





| Report of Analysis | | | | | | | | |
|---|------------------------|--|-------------------------------|-----------|----------------|------------|------------------------------|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC SC SV | W-9 (19.4') 22892-1) - Soil V846 8260C 02-2024 Crop | SW846 5035 sey Avenue, Bro | ooklyn, N | Date Perce | I | 06/22/16 06/24/16 76.1 | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| Run #1 Run #2 | Y164284.E |) 1 | 06/27/16 | PS | 06/24/16 10:00 | n/a | VY7164 | |
| | Initial Wei | ght | | | | | | |
| Run #1 Run #2 | 5.3 g | | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 7.7 | 12 | 2.3 | ug/kg | J |
| 71-43-2 | Benzene | ND | 0.62 | 0.15 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.2 | 0.40 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.5 | 0.19 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.2 | 0.33 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.2 | 0.60 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 12 | 2.2 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.5 | 0.21 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.5 | 0.21 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.5 | 0.20 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.2 | 0.53 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.5 | 0.30 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.2 | 0.26 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.5 | 0.68 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.5 | 0.60 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.5 | 0.19 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.2 | 0.30 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.2 | 0.21 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.2 | 0.17 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.2 | 0.19 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.2 | 0.68 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.2 | 0.23 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.2 | 0.21 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.2 | 0.19 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.2 | 0.54 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.2 | 0.20 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.5 | 0.38 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.5 | 0.24 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.5 | 0.27 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.2 | 0.18 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.2 | 0.60 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.2 | 1.7 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | MW-9 (19.4') JC22892-1 SO - Soil SW846 8260C SW84 2002-2024 Cropsey A | | klyn NY | | Date | Sampled: Received: nt Solids: | 06/22/16 06/24/16 76.1 |
|---|---|--------|---------|-----|-------|-------------------------------------|------------------------------|
| VOA TCL List | pound | Result | RL | MDL | Units | 0 | |

| | 1 | | | | | • |
|------------|----------------------------|--------|--------|-----------------------|----------|---|
| 98-82-8 | Isopropylbenzene | ND | 2.5 | 0.19 | ug/kg | |
| 79-20-9 | Methyl Acetate | ND | 6.2 | 2.5 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 2.5 | 0.63 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.2 | 0.33 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.2 | 1.1 | ug/kg | |
| 75-09-2 | Methylene chloride | 1.5 | 6.2 | 0.42 | ug/kg | J |
| 100-42-5 | Styrene | ND | 2.5 | 0.18 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.5 | 0.30 | ug/kg | |
| 127-18-4 | Tetrachloroethene | ND | 2.5 | 0.35 | ug/kg | |
| 108-88-3 | Toluene | ND | 1.2 | 0.15 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.2 | 0.28 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.2 | 0.23 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.5 | 0.21 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.5 | 0.40 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 1.2 | 0.24 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 6.2 | 0.78 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 2.5 | 0.25 | ug/kg | |
| | m,p-Xylene | ND | 1.2 | 0.27 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 1.2 | 0.25 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 1.2 | 0.25 | ug/kg | |
| | | | | | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limit | ts | |
| 10.00 50 5 | | 0.004 | | 5 0 4 0 | . | |
| 1868-53-7 | Dibromofluoromethane | 98% | | 70-12 | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 90% | | 68-12 | | |
| 2037-26-5 | Toluene-D8 | 96% | | 77-12 | | |
| 460-00-4 | 4-Bromofluorobenzene | 100% | | 72-13 | 0% | |
| | | | | | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





Page 2 of 2

| | | | nope | | 11141.5515 | | ruge r or r | |
|-----------|---------------|------------------------|-----------------|-----------|----------------|--------------------------------|------------------|--|
| Client Sa | mple ID: MW- | 9 (30') | | | | | | |
| Lab Sam | ple ID: JC22 | 892-2 | | | Date | Date Sampled: 06 | | |
| Matrix: | SO - | Soil | | | Date | Date Received: 06/24/16 | | |
| Method: | SW84 | SW846 8260C SW846 5035 | | | Perc | 5.1 | | |
| Project: | 2002- | 2024 Crop | sey Avenue, Bro | ooklyn, N | IΥ | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| Run #1 | Y164251.D | 1 | 06/25/16 | PS | 06/24/16 10:00 | - | VY7163 | |
| Run #2 | | | | | | | | |
| | Initial Weigh | t | | | | | | |
| Run #1 | 5.4 g | | | | | | | |
| Run #2 | - | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 11 | 2.0 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.54 | 0.13 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.4 | 0.35 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.2 | 0.17 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.4 | 0.29 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.4 | 0.53 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 11 | 1.9 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.2 | 0.18 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.2 | 0.18 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.2 | 0.18 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.4 | 0.47 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.2 | 0.26 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.4 | 0.23 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.2 | 0.59 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.2 | 0.53 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.2 | 0.16 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.1 | 0.26 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.1 | 0.19 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.1 | 0.15 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.1 | 0.17 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.4 | 0.59 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.1 | 0.20 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.1 | 0.19 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.1 | 0.17 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | 0.82 | 1.1 | 0.48 | ug/kg | J |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.1 | 0.17 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.2 | 0.34 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.2 | 0.21 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.2 | 0.24 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.1 | 0.16 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.4 | 0.53 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.4 | 1.5 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





11 of 404

E = Indicates value exceeds calibration range

79-01-6

75-69-4

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

Trichloroethene

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

o-Xylene

Trichlorofluoromethane

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Report of Analysis

| Client Sam Lab Sample Matrix: Method: Project: | e ID: JC22892-2 SO - Soil SW846 8260C SW846 2002-2024 Cropsey Av | | klyn, NY | | Date | Sampled: Received: ent Solids: | 06/22/16 06/24/16 85.1 | |
|--|---|--------|----------|------|-------|--------------------------------------|------------------------------|--|
| VOA TCL | List | | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | | |
| 98-82-8 | Isopropylbenzene | ND | 2.2 | 0.17 | ug/kg | | | |
| 79-20-9 | Methyl Acetate | ND | 5.4 | 2.2 | ug/kg | | | |
| 108-87-2 | Methylcyclohexane | ND | 2.2 | 0.55 | ug/kg | | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.1 | 0.29 | ug/kg | | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.4 | 0.92 | ug/kg | | | |
| 75-09-2 | Methylene chloride | 1.2 | 5.4 | 0.37 | ug/kg | J | | |
| 100-42-5 | Styrene | ND | 2.2 | 0.16 | ug/kg | | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.2 | 0.26 | ug/kg | | | |
| 127-18-4 | Tetrachloroethene | 2.8 | 2.2 | 0.31 | ug/kg | | | |
| 108-88-3 | Toluene | ND | 1.1 | 0.14 | ug/kg | | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.4 | 0.25 | ug/kg | | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.4 | 0.20 | ug/kg | | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.2 | 0.18 | ug/kg | | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.2 | 0.35 | ug/kg | | | |

1.1

5.4

2.2

1.1

1.1

1.1

Run# 2

0.21

0.68

0.22

0.24

0.22

0.22

Limits

70-122%

68-124%

77-125%

72-130%

ND

ND

ND

ND

ND

ND

Run#1

99%

96%

99%

99%

ND = Not detectedMDL = Method Detection Limit RL = Reporting Limit E = Indicates value exceeds calibration range

J = Indicates an estimated value

ug/kg

ug/kg

ug/kg

ug/kg

ug/kg

ug/kg

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2

SGS

12 of 404

| Report of Analysis Pag | | | | | | | | | | | |
|---|----------------------------|--|-------------------------------|-----------------|----|---------------------------------|--------------------------|------------------------------|--|--|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: J S S | 4W-10 (19.5') C22892-3 O - Soil W846 8260C 002-2024 Crop | SW846 5035 sey Avenue, Bro | ooklyn, | NY | Date | Received: | 06/22/16 06/24/16 78.0 | | | |
| Run #1 Run #2 | File ID Y164285. | DF D 1 | Analyzed 06/27/16 | By PS | | Prep Date 06/24/16 11:00 | Prep Batch n/a | n Analytical Batch VY7164 | | | |
| Run #1 Run #2 | Initial W 4.4 g | eight | | | | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 15 | 2.7 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.73 | 0.17 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 7.3 | 0.46 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.9 | 0.22 | ug/kg | |
| 75-25-2 | Bromoform | ND | 7.3 | 0.39 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 7.3 | 0.71 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 15 | 2.6 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.9 | 0.25 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.9 | 0.24 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.9 | 0.24 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 7.3 | 0.63 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.9 | 0.35 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 7.3 | 0.31 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.9 | 0.80 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.9 | 0.71 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.9 | 0.22 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.5 | 0.35 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.5 | 0.25 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.5 | 0.20 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.5 | 0.22 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 7.3 | 0.79 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.5 | 0.27 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.5 | 0.25 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.5 | 0.22 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.5 | 0.64 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.5 | 0.23 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.9 | 0.45 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.9 | 0.29 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.9 | 0.32 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.5 | 0.22 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 7.3 | 0.71 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 7.3 | 2.0 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



JC22892

4.3 4

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Compound

CAS No.

Report of Analysis

| | 2892-3 | Date Sampled: | 06/00/16 |
|---------------|-------------------------------------|-----------------|----------|
| _ | | Date Sampleu. | 06/22/16 |
| Matrix: SO | - Soil | Date Received: | 06/24/16 |
| Method: SW8 | 846 8260C SW846 5035 | Percent Solids: | 78.0 |
| Project: 2002 | 2-2024 Cropsey Avenue, Brooklyn, NY | | |

RL

MDL

Units

Q

Result

| C/10 110. | Compound | Result | KL | MIDL | Cints | Y |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 2.9 | 0.22 | ug/kg | |
| 79-20-9 | Methyl Acetate | ND | 7.3 | 3.0 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 2.9 | 0.74 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.5 | 0.39 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 7.3 | 1.2 | ug/kg | |
| 75-09-2 | Methylene chloride | 1.2 | 7.3 | 0.50 | ug/kg | J |
| 100-42-5 | Styrene | ND | 2.9 | 0.21 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.9 | 0.35 | ug/kg | |
| 127-18-4 | Tetrachloroethene | 0.77 | 2.9 | 0.41 | ug/kg | J |
| 108-88-3 | Toluene | ND | 1.5 | 0.18 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 7.3 | 0.33 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 7.3 | 0.27 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.9 | 0.24 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.9 | 0.47 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 1.5 | 0.28 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 7.3 | 0.92 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 2.9 | 0.29 | ug/kg | |
| | m,p-Xylene | ND | 1.5 | 0.32 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 1.5 | 0.29 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 1.5 | 0.29 | ug/kg | |
| | | | | | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1060 52 7 | | 000/ | | 70.10 | 20/ | |
| 1868-53-7 | Dibromofluoromethane | 99% | | 70-12 | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 91% | | 68-12 | | |
| 2037-26-5 | Toluene-D8 | 97% | | 77-12 | | |
| 460-00-4 | 4-Bromofluorobenzene | 99% | | 72-13 | 0% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2



| | | | nope | | u iy 515 | | 1 ugo 1 01 2 |
|-----------|--------------|--|----------|----|-----------------|------------|------------------|
| Client Sa | mple ID: MV | V-10 (30') | | | | | |
| Lab Sam | ple ID: JC2 | 22892-4 | | | Date | Sampled: 0 | 6/22/16 |
| Matrix: | SO | - Soil | | | Date | 6/24/16 | |
| Method: | SW | SW846 8260C SW846 5035 | | | Perc | 9.3 | |
| Project: | 200 | 2002-2024 Cropsey Avenue, Brooklyn, NY | | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | Y164263.D | 1 | 06/25/16 | PS | 06/24/16 11:00 | n/a | VY7163 |
| Run #2 | | | | | | | |
| | Initial Weig | ght | | | | | |
| Run #1 | 4.8 g | | | | | | |
| Run #2 | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 13 | 2.5 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.66 | 0.16 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.6 | 0.42 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.6 | 0.20 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.6 | 0.35 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.6 | 0.64 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 13 | 2.3 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.6 | 0.22 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.6 | 0.22 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.6 | 0.21 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.6 | 0.56 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.6 | 0.31 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.6 | 0.28 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.6 | 0.72 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.6 | 0.64 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.6 | 0.20 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.3 | 0.32 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.3 | 0.22 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.3 | 0.18 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.3 | 0.20 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.6 | 0.72 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.3 | 0.25 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.3 | 0.22 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.3 | 0.20 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.3 | 0.58 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.3 | 0.21 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.6 | 0.41 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.6 | 0.26 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.6 | 0.29 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.3 | 0.20 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.6 | 0.64 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.6 | 1.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

4-Methyl-2-pentanone(MIBK)

1,1,2,2-Tetrachloroethane

1,2,3-Trichlorobenzene

1,2,4-Trichlorobenzene

1,1,1-Trichloroethane

1,1,2-Trichloroethane

Trichlorofluoromethane

Surrogate Recoveries

Dibromofluoromethane

1.2-Dichloroethane-D4

4-Bromofluorobenzene

Trichloroethene

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

o-Xylene

Methylene chloride

Tetrachloroethene

Styrene

Toluene

108-10-1

100-42-5

79-34-5

127-18-4

108-88-3

87-61-6

120-82-1

71-55-6

79-00-5

79-01-6

75-69-4

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

75-09-2

Report of Analysis

| Client Sam Lab Sampl Matrix: Method: Project: | 1 () | | oklyn, NY | | Date | Sampled: Received: ent Solids: | 06/22/16 06/24/16 79.3 |
|---|-------------------------|--------|-----------|------|-------|--------------------------------------|------------------------------|
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 2.6 | 0.20 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 6.6 | 2.7 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 2.6 | 0.66 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.3 | 0.35 | ug/kg | | |

6.6

6.6

2.6

2.6

2.6

1.3

6.6

6.6

2.6

2.6

1.3

6.6

2.6

1.3

1.3

1.3

Run# 2

1.1

0.45

0.19

0.31

0.37

0.16

0.30

0.24

0.22

0.42

0.25

0.83

0.27

0.29

0.27

0.27

Limits

70-122%

68-124%

77-125%

72-130%

ug/kg

J

J

ND

1.2

ND

ND

1.1

ND

99%

94%

99%

99%

Run#1

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit E = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



16 of 404

IC22892

Page 2 of 2

| Report of Analysis Pa | | | | | | | | | | |
|---|----------------------------|--|--------------------------|-----------------|---------------------------------|--------------------------|-----------------------------------|--|--|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: J S S | 4W-8 (19.5') C22892-5 O - Soil W846 8260C 002-2024 Crops | | ooklyn, N | Date Perc | I I | 06/22/16 06/24/16 79.8 | | | |
| Run #1 Run #2 | File ID Y164257. | DF D 1 | Analyzed 06/25/16 | By PS | Prep Date 06/24/16 11:00 | Prep Batch n/a | Analytical Batch VY7163 | | | |
| Run #1 Run #2 | Initial W 5.0 g | eight | | | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 13 | 2.3 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.63 | 0.15 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.3 | 0.40 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.5 | 0.19 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.3 | 0.33 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.3 | 0.61 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 13 | 2.2 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.5 | 0.21 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.5 | 0.21 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.5 | 0.20 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.3 | 0.54 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.5 | 0.30 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.3 | 0.26 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.5 | 0.68 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.5 | 0.61 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.5 | 0.19 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.3 | 0.30 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.3 | 0.21 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.3 | 0.17 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.3 | 0.19 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.3 | 0.68 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.3 | 0.23 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.3 | 0.21 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.3 | 0.19 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.3 | 0.55 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.3 | 0.20 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.5 | 0.39 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.5 | 0.25 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.5 | 0.28 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.3 | 0.19 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.3 | 0.61 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.3 | 1.7 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

 $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Sample II Lab Sample ID: Matrix: Method: Project: | D: MW-8 (19.5') JC22892-5 SO - Soil SW846 8260C SW 2002-2024 Cropsey | | klyn, NY | | Date | Sampled: Received: ent Solids: | |
|--|--|--------|----------|-----|-------|--------------------------------------|--|
| VOA TCL List | | | | | | | |
| CAS No. Con | npound | Result | RL | MDL | Units | Q | |

| | 1 | | | | | |
|------------|-------------------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 2.5 | 0.19 | ug/kg | |
| 79-20-9 | Methyl Acetate | ND | 6.3 | 2.5 | ug/kg | |
| 108-87-2 | Methylcyclohexane | ND | 2.5 | 0.63 | ug/kg | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.3 | 0.33 | ug/kg | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.3 | 1.1 | ug/kg | |
| 75-09-2 | Methylene chloride | 1.2 | 6.3 | 0.43 | ug/kg | J |
| 100-42-5 | Styrene | ND | 2.5 | 0.18 | ug/kg | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.5 | 0.30 | ug/kg | |
| 127-18-4 | Tetrachloroethene | ND | 2.5 | 0.35 | ug/kg | |
| 108-88-3 | Toluene | ND | 1.3 | 0.16 | ug/kg | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.3 | 0.28 | ug/kg | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.3 | 0.23 | ug/kg | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.5 | 0.21 | ug/kg | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.5 | 0.40 | ug/kg | |
| 79-01-6 | Trichloroethene | ND | 1.3 | 0.24 | ug/kg | |
| 75-69-4 | Trichlorofluoromethane | ND | 6.3 | 0.79 | ug/kg | |
| 75-01-4 | Vinyl chloride | ND | 2.5 | 0.25 | ug/kg | |
| | m,p-Xylene | ND | 1.3 | 0.27 | ug/kg | |
| 95-47-6 | o-Xylene | ND | 1.3 | 0.25 | ug/kg | |
| 1330-20-7 | Xylene (total) | ND | 1.3 | 0.25 | ug/kg | |
| | | | | | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 100% | | 70.12 | 20/ | |
| | | | | 70-12 | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 Toluene-D8 | 95% | | 68-12 | | |
| 2037-26-5 | | 100% | | 77-12 | | |
| 460-00-4 | 4-Bromofluorobenzene | 100% | | 72-13 | 0% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



18 of 404

JC22892

Page 2 of 2

4.5 **4**

| | | | | | 1141 9 515 | 1 age 1 01 2 | |
|-----------|--------------|-------------|-----------------|------------|----------------|---------------|------------------|
| Client Sa | mple ID: MW | (-8 (30') | | | | | |
| Lab Sam | ple ID: JC2 | 2892-6 | | | Date | Sampled: (| 06/22/16 |
| Matrix: | SO - | Soil | | | Date | Received: (| 06/24/16 |
| Method: | SW8 | 846 8260C | SW846 5035 | | Perc | ent Solids: 8 | 81.1 |
| Project: | 2002 | 2-2024 Crop | sey Avenue, Bro | ooklyn, NY | 7 | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | Y164286.D | 1 | 06/27/16 | PS | 06/24/16 11:00 | n/a | VY7164 |
| Run #2 | | | | | | | |
| | Initial Weig | ht | | | | | |
| Run #1 | 5.5 g | | | | | | |
| Run #2 | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | 29.6 | 11 | 2.1 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.56 | 0.13 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.6 | 0.36 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.2 | 0.17 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.6 | 0.30 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.6 | 0.54 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 11 | 2.0 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.2 | 0.19 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.2 | 0.19 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.2 | 0.18 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.6 | 0.48 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.2 | 0.27 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.6 | 0.24 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.2 | 0.61 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.2 | 0.54 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.2 | 0.17 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.1 | 0.27 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.1 | 0.19 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.1 | 0.15 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.1 | 0.17 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.6 | 0.61 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.1 | 0.21 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.1 | 0.19 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.1 | 0.17 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.1 | 0.49 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.1 | 0.18 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.2 | 0.35 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.2 | 0.22 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.2 | 0.25 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.1 | 0.17 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.6 | 0.54 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.6 | 1.6 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound

Page 1 of 2

19 of 404

E = Indicates value exceeds calibration range

J = Indicates an estimated value

79-01-6

75-69-4

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

Trichloroethene

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

o-Xylene

Trichlorofluoromethane

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Report of Analysis

| Client Samj Lab Sample Matrix: Method: Project: | | | Date | Sampled: Received: ent Solids: | 06/22/16 06/24/16 81.1 | | |
|---|----------------------------|--------|------|--------------------------------------|------------------------------|---|--|
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 2.2 | 0.17 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 5.6 | 2.3 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 2.2 | 0.57 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.1 | 0.30 | ug/kg | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.6 | 0.95 | ug/kg | | |
| 75-09-2 | Methylene chloride | 0.51 | 5.6 | 0.38 | ug/kg | J | |
| 100-42-5 | Styrene | ND | 2.2 | 0.16 | ug/kg | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.2 | 0.27 | ug/kg | | |
| 127-18-4 | Tetrachloroethene | 0.48 | 2.2 | 0.31 | ug/kg | J | |
| 108-88-3 | Toluene | ND | 1.1 | 0.14 | ug/kg | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.6 | 0.25 | ug/kg | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.6 | 0.20 | ug/kg | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.2 | 0.19 | ug/kg | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.2 | 0.36 | ug/kg | | |

1.1

5.6

2.2

1.1

1.1

1.1

Run# 2

0.21

0.71

0.23

0.25

0.23

0.23

Limits

70-122%

68-124%

77-125%

72-130%

ug/kg

ug/kg

ug/kg

ug/kg

ug/kg

ug/kg

ND

ND

ND

ND

ND

ND

99%

91%

97%

101%

Run#1

| ND = Not detected | MDL = Method Detection Limit |
|--------------------------|------------------------------|
| RL = Reporting Limit | |
| E = Indicates value exce | eeds calibration range |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



SGS

4.6



| Client San Lab Samp Matrix: Method: Project: | ole ID: JC228 SO - S SW84 | Soil 6 8260C S | W846 5035 y Avenue, Br | ooklyn, NY | | Date | Received: 06 | 5/23/16 5/24/16 0.6 |
|--|---------------------------------|-------------------|---------------------------|------------|---------|--------------------|--------------|---------------------------|
| | File ID | DF | Analyzed | By | Prep D | ate | Prep Batch | Analytical Batch |
| Run #1 | Y164280.D | 1 | 06/27/16 | PS | 06/24/1 | 16 11:00 | n/a | VY7164 |
| Run #2 | E236298.D | 1 | 06/29/16 | TDN | 06/24/1 | 16 11:00 | n/a | VE10266 |
| | Initial Weight | Final Vo | olume | Methanol A | liquot | | | |
| Run #1 | 5.2 g | | | | | | | |
| Run #2 | 4.5 g | 5.0 ml | | 100 ul | | | | |
| VOA TCI | List | | | | | | | |
| CAS No. | Compound | | Result | RL | MDL | Units | Q | |
| 67-64-1 | Acetone | | ND | 12 | 2.2 | ug/kg | | |
| 71-43-2 | Benzene | | ND | 0.60 | 0.14 | ug/kg | | |
| 74-97-5 | Bromochloro | methane | ND | 6.0 | 0.38 | ug/kg | | |
| 75 27 1 | Promodiable | omothono | ND | 2.4 | 0.10 | $n\alpha/lc\alpha$ | | |

Report of Analysis

| | | | | | 00 |
|------------|-----------------------------|------|-----|------|-------|
| 74-97-5 | Bromochloromethane | ND | 6.0 | 0.38 | ug/kg |
| 75-27-4 | Bromodichloromethane | ND | 2.4 | 0.18 | ug/kg |
| 75-25-2 | Bromoform | ND | 6.0 | 0.32 | ug/kg |
| 74-83-9 | Bromomethane | ND | 6.0 | 0.58 | ug/kg |
| 78-93-3 | 2-Butanone (MEK) | ND | 12 | 2.1 | ug/kg |
| 75-15-0 | Carbon disulfide | ND | 2.4 | 0.20 | ug/kg |
| 56-23-5 | Carbon tetrachloride | ND | 2.4 | 0.20 | ug/kg |
| 108-90-7 | Chlorobenzene | ND | 2.4 | 0.19 | ug/kg |
| 75-00-3 | Chloroethane | ND | 6.0 | 0.51 | ug/kg |
| 67-66-3 | Chloroform | ND | 2.4 | 0.28 | ug/kg |
| 74-87-3 | Chloromethane | ND | 6.0 | 0.25 | ug/kg |
| 110-82-7 | Cyclohexane | ND | 2.4 | 0.65 | ug/kg |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.4 | 0.58 | ug/kg |
| 124-48-1 | Dibromochloromethane | ND | 2.4 | 0.18 | ug/kg |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.2 | 0.29 | ug/kg |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.2 | 0.20 | ug/kg |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.2 | 0.16 | ug/kg |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.2 | 0.18 | ug/kg |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.0 | 0.65 | ug/kg |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.2 | 0.22 | ug/kg |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.2 | 0.20 | ug/kg |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.2 | 0.18 | ug/kg |
| 156-59-2 | cis-1,2-Dichloroethene | 30.8 | 1.2 | 0.52 | ug/kg |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.2 | 0.19 | ug/kg |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.4 | 0.37 | ug/kg |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.4 | 0.23 | ug/kg |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.4 | 0.26 | ug/kg |
| 100-41-4 | Ethylbenzene | ND | 1.2 | 0.18 | ug/kg |
| 76-13-1 | Freon 113 | ND | 6.0 | 0.58 | ug/kg |
| 591-78-6 | 2-Hexanone | ND | 6.0 | 1.7 | ug/kg |
| | | | | | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 1 of 2

21 of 404

| Client Sample ID: MW-51 (20') Lab Sample ID: JC22892-7 Matrix: SO - Soil Method: SW846 8260C SW84 Project: 2002-2024 Cropsey A | | | klyn, NY | | Date | Sampled: Received: ent Solids: | 06/23/16 06/24/16 80.6 |
|--|------------------------------|------------------|----------|------|-------|--------------------------------------|------------------------------|
| VOA TCL | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 98-82-8 | Isopropylbenzene | ND | 2.4 | 0.18 | ug/kg | | |
| 79-20-9 | Methyl Acetate | ND | 6.0 | 2.4 | ug/kg | | |
| 108-87-2 | Methylcyclohexane | ND | 2.4 | 0.60 | ug/kg | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.2 | 0.32 | ug/kg | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.0 | 1.0 | ug/kg | | |
| 75-09-2 | Methylene chloride | 1.4 | 6.0 | 0.41 | ug/kg | J | |
| 100-42-5 | Styrene | ND | 2.4 | 0.17 | ug/kg | | |
| 79-34-5 | 1, 1, 2, 2-Tetrachloroethane | ND | 2.4 | 0.29 | ug/kg | | |
| 127-18-4 | Tetrachloroethene | 736 ^a | 160 | 23 | ug/kg | | |
| 108-88-3 | Toluene | ND | 1.2 | 0.15 | ug/kg | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.0 | 0.27 | ug/kg | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.0 | 0.22 | ug/kg | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.4 | 0.20 | ug/kg | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.4 | 0.39 | ug/kg | | |
| 79-01-6 | Trichloroethene | 3.9 | 1.2 | 0.23 | ug/kg | | |
| 75-69-4 | Trichlorofluoromethane | ND | 6.0 | 0.75 | ug/kg | | |

2.4

1.2

1.2

1.2

Run# 2

103%

94%

98%

98%

0.24

0.26

0.24

0.24

Limits

70-122%

68-124%

77-125%

72-130%

ND

ND

ND

ND

Run#1

98%

91%

98%

99%

(a) Result is from Run# 2

Vinyl chloride

m,p-Xylene

Xylene (total)

Toluene-D8

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

o-Xylene

75-01-4

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

J = Indicates an estimated value

ug/kg

ug/kg

ug/kg

ug/kg

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| | | | nope | | 1413.515 | | ruge rorz | | |
|-----------|--------------------------------|-------------|-----------------|------------|----------------------|--------------|------------------|--|--|
| Client Sa | mple ID: MW | -5I (23') | | | | | | | |
| Lab Sam | ple ID: JC22 | 2892-8 | | | Date | Sampled: 06 | 5/23/16 | | |
| Matrix: | SO - | Soil | | | Date | Received: 06 | 5/24/16 | | |
| Method: | Method: SW846 8260C SW846 5035 | | | | Percent Solids: 78.0 | | | | |
| Project: | 2002 | 2-2024 Crop | sey Avenue, Bro | ooklyn, NY | | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | | |
| Run #1 | Y164287.D | 1 | 06/27/16 | PS | 06/24/16 11:00 | n/a | VY7164 | | |
| Run #2 | | | | | | | | | |
| | Initial Weig | nt | | | | | | | |
| Run #1 | 5.4 g | | | | | | | | |
| Run #2 | | | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 12 | 2.2 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.59 | 0.14 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 5.9 | 0.38 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.4 | 0.18 | ug/kg | |
| 75-25-2 | Bromoform | ND | 5.9 | 0.32 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 5.9 | 0.58 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 12 | 2.1 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.4 | 0.20 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.4 | 0.20 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.4 | 0.19 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 5.9 | 0.51 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.4 | 0.28 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 5.9 | 0.25 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.4 | 0.65 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.4 | 0.57 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.4 | 0.18 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.2 | 0.29 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.2 | 0.20 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.2 | 0.16 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.2 | 0.18 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 5.9 | 0.65 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.2 | 0.22 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.2 | 0.20 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.2 | 0.18 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.2 | 0.52 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.2 | 0.19 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.4 | 0.37 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.4 | 0.23 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.4 | 0.26 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.2 | 0.18 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 5.9 | 0.57 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 5.9 | 1.7 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

23 of 404

JC22892

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Samp Lab Sample Matrix: Method: Project: | | | | | Date | Sampled: Received: ent Solids: | 06/23/16 06/24/16 78.0 | | |
|---|----------------------------|--------|-----|------|-------|--------------------------------------|------------------------------|--|--|
| VOA TCL I | List | | | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | | | |
| 98-82-8 | Isopropylbenzene | ND | 2.4 | 0.18 | ug/kg | | | | |
| 79-20-9 | Methyl Acetate | ND | 5.9 | 2.4 | ug/kg | | | | |
| 108-87-2 | Methylcyclohexane | ND | 2.4 | 0.60 | ug/kg | | | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.2 | 0.31 | ug/kg | | | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.9 | 1.0 | ug/kg | | | | |
| 75-09-2 | Methylene chloride | 1.1 | 5.9 | 0.41 | ug/kg | J | | | |
| 100-42-5 | Styrene | ND | 2.4 | 0.17 | ug/kg | | | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.4 | 0.28 | ug/kg | | | | |
| 127-18-4 | Tetrachloroethene | 3.6 | 2.4 | 0.33 | ug/kg | | | | |
| 108-88-3 | Toluene | ND | 1.2 | 0.15 | ug/kg | | | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 5.9 | 0.27 | ug/kg | | | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.9 | 0.22 | ug/kg | | | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.4 | 0.20 | ug/kg | | | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.4 | 0.38 | ug/kg | | | | |
| 79-01-6 | Trichloroethene | 0.26 | 1.2 | 0.23 | ug/kg | J | | | |
| 75-69-4 | Trichlorofluoromethane | ND | 5.9 | 0.75 | ug/kg | | | | |
| 75-01-4 | Vinyl chloride | ND | 2.4 | 0.24 | ug/kg | | | | |
| | m,p-Xylene | ND | 1.2 | 0.26 | ug/kg | | | | |
| | | | | | | | | | |

1.2

1.2

Run# 2

0.24

0.24

Limits

70-122%

68-124%

77-125%

72-130%

ND

ND

Run#1

97%

90%

97%

100%

95-47-6

1330-20-7

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

o-Xylene

Xylene (total)

Toluene-D8

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

J = Indicates an estimated value

ug/kg

ug/kg

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4.8

Page 2 of 2



| | Report of Analysis Pa | | | | | | | | | | |
|---|------------------------------|---|--------------------------|-----------------|----------------------|---------------------------------|-------------------------|------------------------------|--|--|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: | MW-5I (31') JC22892-9 SO - Soil SW846 8260C SW846 5035 2002-2024 Cropsey Avenue, Brooklyn, NY | | | Date Date Perc | 06/23/16 06/24/16 78.7 | | | | | |
| Run #1 Run #2 | File ID Y16426 | DF 1.D 1 | Analyzed 06/25/16 | By PS | | Prep Date 06/24/16 11:00 | Prep Batc n/a | h Analytical Batch VY7163 | | | |
| Run #1 Run #2 | Initial V 4.9 g | Veight | | | | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 13 | 2.4 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.65 | 0.16 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.5 | 0.41 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.6 | 0.20 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.5 | 0.34 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.5 | 0.63 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 13 | 2.3 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.6 | 0.22 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.6 | 0.22 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.6 | 0.21 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.5 | 0.56 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.6 | 0.31 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.5 | 0.27 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.6 | 0.71 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.6 | 0.63 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.6 | 0.19 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.3 | 0.31 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.3 | 0.22 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.3 | 0.18 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.3 | 0.20 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.5 | 0.71 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.3 | 0.24 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.3 | 0.22 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.3 | 0.20 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.3 | 0.57 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.3 | 0.20 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.6 | 0.40 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.6 | 0.25 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.6 | 0.29 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.3 | 0.19 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.5 | 0.63 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.5 | 1.8 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



25 of 404

ACCUTEST JC22892



E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Sam Lab Sample Matrix: Method: Project: | e ID: JC22892-9 SO - Soil SW846 8260C SW846 | JC22892-9 | | | | | 06/23/16 06/24/16 78.7 | | | | |
|--|---|-----------|-----|------|-------|---|------------------------------|--|--|--|--|
| VOA TCL List | | | | | | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | | | | | |
| 98-82-8 | Isopropylbenzene | ND | 2.6 | 0.20 | ug/kg | | | | | | |
| 79-20-9 | Methyl Acetate | ND | 6.5 | 2.6 | ug/kg | | | | | | |
| 108-87-2 | Methylcyclohexane | ND | 2.6 | 0.65 | ug/kg | | | | | | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.3 | 0.34 | ug/kg | | | | | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 6.5 | 1.1 | ug/kg | | | | | | |
| 75-09-2 | Methylene chloride | 0.98 | 6.5 | 0.44 | ug/kg | J | | | | | |
| 100-42-5 | Styrene | ND | 2.6 | 0.19 | ug/kg | | | | | | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 2.6 | 0.31 | ug/kg | | | | | | |
| 127-18-4 | Tetrachloroethene | 2.2 | 2.6 | 0.36 | ug/kg | J | | | | | |
| 108-88-3 | Toluene | ND | 1.3 | 0.16 | ug/kg | | | | | | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 6.5 | 0.29 | ug/kg | | | | | | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 6.5 | 0.24 | ug/kg | | | | | | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 2.6 | 0.22 | ug/kg | | | | | | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 2.6 | 0.42 | ug/kg | | | | | | |
| 79-01-6 | Trichloroethene | ND | 1.3 | 0.25 | ug/kg | | | | | | |
| 75-69-4 | Trichlorofluoromethane | ND | 6.5 | 0.82 | ug/kg | | | | | | |
| 75-01-4 | Vinyl chloride | ND | 2.6 | 0.26 | ug/kg | | | | | | |
| | m,p-Xylene | ND | 1.3 | 0.28 | ug/kg | | | | | | |
| 95-47-6 | o-Xylene | ND | 1.3 | 0.26 | ug/kg | | | | | | |
| 1330-20-7 | Xylene (total) | ND | 1.3 | 0.26 | ug/kg | | | | | | |

Run#1

99%

94%

99%

98%

Run# 2

Limits

70-122%

68-124%

77-125%

72-130%

| ND = Not detected | MDL = Method Detection Limit |
|--------------------------|------------------------------|
| RL = Reporting Limit | |
| E = Indicates value exce | eeds calibration range |

Surrogate Recoveries

Dibromofluoromethane

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Toluene-D8

CAS No.

1868-53-7

2037-26-5

460-00-4

17060-07-0

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound

Page 2 of 2



| | Report of Analysis Pa | | | | | | | | | | |
|---|------------------------------|--|----|--------------------------|----------------------|------------------------------|---------------------------------|-------------------------|------------------------------|--|--|
| Client San Lab Sam Matrix: Method: Project: | - | MW-5I (50') JC22892-10 SO - Soil SW846 8260C SW846 5035 2002-2024 Cropsey Avenue, Brooklyn, NY | | | Date Date Perc | 06/23/16 06/24/16 77.3 | | | | | |
| Run #1 Run #2 | File ID Y16426 | - | DF | Analyzed 06/25/16 | By PS | | Prep Date 06/24/16 11:00 | Prep Batc n/a | h Analytical Batch VY7163 | | |
| Run #1 Run #2 | Initial 4.7 g | Veight | | | | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 14 | 2.6 | ug/kg | |
| 71-43-2 | Benzene | ND | 0.69 | 0.17 | ug/kg | |
| 74-97-5 | Bromochloromethane | ND | 6.9 | 0.44 | ug/kg | |
| 75-27-4 | Bromodichloromethane | ND | 2.8 | 0.21 | ug/kg | |
| 75-25-2 | Bromoform | ND | 6.9 | 0.37 | ug/kg | |
| 74-83-9 | Bromomethane | ND | 6.9 | 0.67 | ug/kg | |
| 78-93-3 | 2-Butanone (MEK) | ND | 14 | 2.4 | ug/kg | |
| 75-15-0 | Carbon disulfide | ND | 2.8 | 0.23 | ug/kg | |
| 56-23-5 | Carbon tetrachloride | ND | 2.8 | 0.23 | ug/kg | |
| 108-90-7 | Chlorobenzene | ND | 2.8 | 0.22 | ug/kg | |
| 75-00-3 | Chloroethane | ND | 6.9 | 0.59 | ug/kg | |
| 67-66-3 | Chloroform | ND | 2.8 | 0.33 | ug/kg | |
| 74-87-3 | Chloromethane | ND | 6.9 | 0.29 | ug/kg | |
| 110-82-7 | Cyclohexane | ND | 2.8 | 0.75 | ug/kg | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.8 | 0.67 | ug/kg | |
| 124-48-1 | Dibromochloromethane | ND | 2.8 | 0.21 | ug/kg | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.4 | 0.33 | ug/kg | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.4 | 0.24 | ug/kg | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.4 | 0.19 | ug/kg | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.4 | 0.21 | ug/kg | |
| 75-71-8 | Dichlorodifluoromethane | ND | 6.9 | 0.75 | ug/kg | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.4 | 0.26 | ug/kg | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.4 | 0.24 | ug/kg | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.4 | 0.21 | ug/kg | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.4 | 0.60 | ug/kg | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.4 | 0.22 | ug/kg | |
| 78-87-5 | 1,2-Dichloropropane | ND | 2.8 | 0.43 | ug/kg | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 2.8 | 0.27 | ug/kg | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 2.8 | 0.30 | ug/kg | |
| 100-41-4 | Ethylbenzene | ND | 1.4 | 0.21 | ug/kg | |
| 76-13-1 | Freon 113 | ND | 6.9 | 0.67 | ug/kg | |
| 591-78-6 | 2-Hexanone | ND | 6.9 | 1.9 | ug/kg | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



27 of 404

E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Lab Sample ID:JC2289Matrix:SO - SoMethod:SW846 | | | 2892-10 | | | | | 06/23/16 06/24/16 77.3 | |
|--|----------------------------|----------------------|---------|--------|---------|-------|---|------------------------------|--|
| VOA TCL | List | | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | | |
| 98-82-8 | Isopro | pylbenzene | ND | 2.8 | 0.21 | ug/kg | | | |
| 79-20-9 | Methy | 1 Acetate | ND | 6.9 | 2.8 | ug/kg | | | |
| 108-87-2 | Methy | lcyclohexane | ND | 2.8 | 0.70 | ug/kg | | | |
| 1634-04-4 | Methy | l Tert Butyl Ether | ND | 1.4 | 0.36 | ug/kg | | | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | | ND | 6.9 | 1.2 | ug/kg | | | |
| 75-09-2 | Methylene chloride | | 1.1 | 6.9 | 0.47 | ug/kg | J | | |
| 100-42-5 | Styren | e | ND | 2.8 | 0.20 | ug/kg | | | |
| 79-34-5 | 1,1,2,1 | 2-Tetrachloroethane | ND | 2.8 | 0.33 | ug/kg | | | |
| 127-18-4 | Tetrac | hloroethene | 2.6 | 2.8 | 0.39 | ug/kg | J | | |
| 108-88-3 | Toluer | ne | ND | 1.4 | 0.17 | ug/kg | | | |
| 87-61-6 | 1,2,3- | Trichlorobenzene | ND | 6.9 | 0.31 | ug/kg | | | |
| 120-82-1 | 1,2,4- | Trichlorobenzene | ND | 6.9 | 0.25 | ug/kg | | | |
| 71-55-6 | 1,1,1- | Trichloroethane | ND | 2.8 | 0.23 | ug/kg | | | |
| 79-00-5 | 1,1,2- | Trichloroethane | ND | 2.8 | 0.44 | ug/kg | | | |
| 79-01-6 | Trichl | oroethene | ND | 1.4 | 0.26 | ug/kg | | | |
| 75-69-4 | Trichl | orofluoromethane | ND | 6.9 | 0.87 | ug/kg | | | |
| 75-01-4 | Vinyl | chloride | ND | 2.8 | 0.28 | ug/kg | | | |
| | m,p-X | ylene | ND | 1.4 | 0.30 | ug/kg | | | |
| 95-47-6 | o-Xyle | ene | ND | 1.4 | 0.28 | ug/kg | | | |
| 1330-20-7 | Xylene | e (total) | ND | 1.4 | 0.28 | ug/kg | | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | Lim | iits | | | |
| 1868-53-7 | Dibroi | Dibromofluoromethane | | | 70-122% | | | | |
| | | | | | | | | | |

94%

99%

99%

1,2-Dichloroethane-D4

4-Bromofluorobenzene

Toluene-D8

17060-07-0

2037-26-5

460-00-4

J = Indicates an estimated value

68-124%

77-125%

72-130%

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2

SGS

| | | | - 1 | | J | | 8 | |
|------------|------------|--|----------|----|-----------|---------------------|------------------|--|
| Client Sam | - | 2002 11 | | | D | | | |
| Lab Samp | le ID: JC2 | 2892-11 | | | Da | Date Sampled: 06/ | | |
| Matrix: | AQ | - Field Bla | unk Soil | | Da | ate Received: (| 6/24/16 | |
| Method: | SW | SW846 8260C | | | Pe | Percent Solids: n/a | | |
| Project: | 200 | 2002-2024 Cropsey Avenue, Brooklyn, NY | | | Y | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| Run #1 | 3A151918.I |) 1 | 06/29/16 | TK | n/a | n/a | V3A6545 | |
| Run #2 | | | | | | | | |
| | Purge Volu | me | | | | | | |
| Run #1 | 5.0 ml | | | | | | | |
| | | | | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.23 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound

Page 1 of 2



E = Indicates value exceeds calibration range

J = Indicates an estimated value

| Client Sample ID: | FB | | |
|-------------------|--|------------------------|----------|
| Lab Sample ID: | JC22892-11 | Date Sampled: | 06/22/16 |
| Matrix: | AQ - Field Blank Soil | Date Received: | 06/24/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | · · · | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|---------------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.22 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 Limits | | its | |
| 1868-53-7 | Dibromofluoromethane | 101% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 103% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 102% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 106% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

SGS

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SGS Accutest LabLink@945686 14:27 10-Feb-2017

| Client Sa Lab Samj Matrix: Method: Project: | ple ID: JC22 AQ - SW8 | BLANK 892-12 Trip Blank 46 8260C -2024 Crop | Soil sey Avenue, Bro | ooklyn, NY | Da | ate Sampled: 06 ate Received: 06 prcent Solids: n/ | |
|---|------------------------------|---|--------------------------|-----------------|-------------------------|--|------------------------------------|
| Run #1 Run #2 | File ID 3A151919.D | DF 1 | Analyzed 06/29/16 | By TK | Prep Date n/a | Prep Batch n/a | Analytical Batch V3A6545 |
| Run #1 Run #2 | Purge Volum 5.0 ml | e | | | | | |

Report of Analysis

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.23 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

 $N = \ Indicates \ presumptive \ evidence \ of \ a \ compound$



Page 1 of 2

SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Client Sample ID: | TRIP BLANK | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC22892-12 | Date Sampled: | 06/23/16 |
| Matrix: | AQ - Trip Blank Soil | Date Received: | 06/24/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|---------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.22 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 100% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 104% | 73-122% | | | |
| 2037-26-5 | Toluene-D8 | 102% | 84-119% | | | |
| 460-00-4 | 4-Bromofluorobenzene | 104% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



32 of 404

JC22892

Appendix F

Soil Boring Logs, Well Logs and Sampling Data Summaries

| | | | - | | | | | _ | | | | |
|----------|-------------------------|---------------------|-------------------------|-----------------|--------|---------------|----------------|--------------------|----------|------------|-------|---|
| | • | | | | | | | | | | ' | Boring #5- MW# Sheet of |
| | | | | | | | | | | | | Project: AMA-Site Investigation |
| | | | | | | | | | | | | JOB # 85265.001 Site: Cropsel AUR |
| | | | | | | | | | | | | Logged By: HRM Proj. Eng: DTS Edited By: |
| | | | | | • | | | | | | | Drilling Contractor: AES |
| | | | | | | | | | | | | Drill RIg Type/Method: 6" Core Orill |
| | | | | | | | | | | | | Drillers Name: Chuck / Tom |
| | | | | | | | | | | | | Borehole Diam./Drill Bit Type Total Depth 4 |
| | | | | | | | | | | | | 1/2"x 2 mario Her. Elev. 100 |
| | | | | | | | | | | | | Hammer Wt: 50 105 Drop: 30" |
| 1 | | | | | | | | | | | | Start Time: OXO Date: 71112 |
| | | | | | | | | | | | | Completion Time: 1100 Date: 711112 |
| | <u>-</u> 5 | Sketch | Мар | of Site | Area V | Vith O | rilling | Loca | tions | | | Backfill Time: Date: |
| | | Ë | | | | Size | | | D | epth | | 1st Water |
| | Sampler Type & Depth | Blaw Counts / 6 In. | (j | Recovered (in.) | | Casing Type & | lier | | j, | | | Boring Depth (ft.) |
| PID/OVA | pler] | Could | Advance (in.) | Verex | | P D | Annulus Filier | | 283 | | | Water Depth (ft.) |
| PID | sam & De | Blow | Adva | Reco | | Casl | Annu | Samolo Doo | Analyses | Feet | | Time %. |
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| r | Boring # S-2 MW# Sheet of |
|--|---|
| | Project: AMA-Site Investigation) |
| | Job # 85265,001 Site: Croosess |
| | Logged By: HPM Proj. Eng: DTS Edited By: |
| · · | Drilling Contractor: AES |
| | Drill Rig Type/Method: 6" Core Drill |
| | Drillers Name: Chuck/Tom |
| | Borehole Dlam./Drill Bit Type Total Depth |
| | 1/2" x 2" MOCYO Ref. Elev. 100 |
| | Hammer Wt: 50105 Drop: 30" |
| | Start Time: 1/00 Date: 7/11/12 |
| | Completion Time: (200 Date: 7/1)/12 |
| Sketch Map of Site Area With Drilling Locations | Backfill Time: Date: |
| C S Depth | 1st Water Borring Depth (ft.) |
| PID/OVA Sampler Type & Depth Advance (in.) Advance (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) | Casing Depth (ft.) |
| PID/OVA Sampler Type & Depth Blow Counts / Advance (in.) Advance (in.) Recovered (in.) Casing Type (Casing Type (Annulus Filler Sample Rec. Analyses | Water Depth (fL) |
| | Time %, |
| 00mc 24"24" | SP, Brass, medium, dry, |
| | med sand |
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|--|---|
| | Boring #5-3 MW# Sheet of |
| | Project: AMA - Site Investigation |
| | Job # 852165,001 Site: Crapsey |
| | Logged By: HRM Proj. Eng: DJS Edited By: |
| | Drilling Contractor: AES |
| | Drill RIg Type/Method: 6 " OOCE Drill |
| | Drillers Name: Chuck / Tom |
| | Borehole Diam./Drill Bit Type Total Depth |
| | 1/a"x a march Ref. Elev. 100 |
| | Hammer Wt: 50165 Drop: 30" |
| | Start Time: 1200 Date: 7/11/12 |
| | Completion Time: 1300 Date: 711112 |
| Sketch Map of Site Area With Drilling Locations | Backfill Time: Date: |
| u u u u u u u u u u u u u u u u u u u | Ist Water Boring Depth (ft.) |
| PID/OVA Sampler Type & Depth & Depth Blow Counts / 6 In. Advance (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered file Annulus Filler Annulus Filler Annulus Filler Feet | Casing Depth (ft.) |
| PID/OVA Sampler Type & Depth & Depth & Depth & Depth & Cashing Type (Annulus Filler Annulus Filler Annulus Filler Annulus Filler Annulus Filler Fiet | Water Depth (fL) |
| PID/OVA Sampler Typ & Depth & Depth Blow Counts Advance (in, Hecovered (in Analyses Analyses Feet | Time 5. |
| p.p.m.e ay ay | SP Brown, Medium, Dry, Med. |
| | sand |
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| HC = Hard Clear | Boring # B-1 MW# Sheet of 2 Project: AMA - Site TDL2SHIGATION Job # 863(65,00) Site: CMDSL Logged By: HPM Proj. Eng: DTS Logged By: HPM Proj. Eng: DTS Logged By: HPM Proj. Eng: DTS Edited By: Drilling Contractor: AES Drilling Contractor: AES Drillers Name: Chuck TOM Borehole Diam./Drill Bit Type I/a " x 5" Hel. Elev. Harnmer Wt: NA Drop: Start Time: IOO Date: 7/13/12 Completion Time: IO3 Date: 7/13/12 |
| Sketch Map of Site Area With Drilling Locations | Backfill Time: Date: |
| PID/OVA Sampler Type & Depth & Depth Blow Counts / 6 In. Advance (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) | 1st Water Boring Depth (ft.) |
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| ٩V ٥ | ipler 3 | v Cou | ance | d (in.) | bu | ulus | de Rec | 103 | | , | | Project: AMA |
| PID/ | Sam Type | Blov | Adva | Ray | Cas | Ann | Same | Analy | Fee | | | Job # 85765,001 Site: CrODSLU |
| | Sampler Type | Blow Counts | | | | | | | 12 13 14 15 16 17 18 19 20 1 2 3 4 5 6 7 8 - 8 - - - - - - - - - - - - - | | | Project: AMA |
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| | · · · · · · · · · · · · · · · · · · · | Boring # B-2 MW# Street (): of 2 Project: AMA - Site Tripstigation Job # 85365, 001 Site: Cropsley Logged By: HPM Proj. Eng: DTS Edited By: Drilling Contractor: AES Drill Rig Type/Method: GEODODE Drillers Name: ODUCK / TOM |
|---|---------------------------------------|--|
| | | Borehole Diam./Drill Bit TypeTotal Depth $1/2$ " \times 5"Ref. Elev. $1/2$ " \times 5"Ref. Elev.Hammer Wt:NADrop:NAStart Time:0900Date:7/13/12Completion Time:0930Date:7/13/12 |
| HC = Hand CRAF Sketch Map of Site Area With Drilling Locatio | 05 | |
| fype nts / 6 in. d (in.) pe & Size | Depth Depth | Backfill Time: Date: 1st Water Boring Depth (ft.) Casing Depth (ft.) Water Depth (ft.) Time Date |
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| r · | Boring # (3~3 MW# Sheet / of a |
| | Project: AMA - Site Incestigation |
| | JOD # 85,765,001 Site: Crop Sel |
| | Logged By: HPM Proj. Eng: DJS Edited By: |
| · | Drilling Contractor: AES |
| | Drill Rig Type/Method: CTEOPIDDE |
| | Drillers Name: Church / Tom |
| | Borehole Diam./Drill Bit Type Total Depth |
| | 1/2" × 5' Ref. Elev. 100 |
| | Hammer WI: NA Drop: NA |
| | Start Time: 1100 Date: 7/13/12 |
| HC = Ham Manr | Completion Time: 1130 Date: 7/13/13 |
| HC = Hand Clear Sketch Map of Site Area With Drilling Locations | Backfill Time: Date: |
| C Depth | 1st Water Boring Depth (ft.) |
| PID/OVA Sampler Type & Depth Blow Counts / 6 In.) Advance (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) Recovered (in.) | Casing Depth (ft.) |
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| | Boring # B-4 MW# Sheet / of 2 |
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| | Project: AMA-Site Investigation |
| | Job # 862/05,00/ Site: (100501) |
| | Logged By: HRM Proj. Eng: DJS Edited By: |
| | Drilling Contractor: AES |
| | Drill Rig Type/Method: Geoprobe |
| | Drillers Name: Chuck / Ton |
| | Borehole Dlam_Drill Bit Type Total Depth |
| | 1/2 ×5' Hef. Elev. 100 |
| | |
| | Start Time: 1030 Date: 7/13/12 |
| HC= Hand Clear | Completion Time: // () Date: 7/13/13 |
| Sketch Map of Site Area With Drilling Locations | Backfill Time: Date: |
| G S Depth | 1st Water Boring Depth (ft.) |
| PID/OVA Sampler Type & Depth & Depth Blow Counts / 6 In.) Advance (in.) Hecovered (in.) Hecovered (in.) Sample Rec. Annulus Filler Annulus Filler Annulus Filler Annulus Filler | Casing Depth (fL) |
| PID/OVA Sampler Type & Depth & Depth Blow Counts / Advance (in.) Advance (in.) Advance (in.) Advance fin.) Sample Rec. Annulus Filler Annulus Filler Annulus Filler Feet | Water Depth (fL) |
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| PID/OVA | Sampler Type | Blow Counts | Advance (In.) | Rcvd (in.) | | Casing | Annulus | Sample Rec. | Analyson | Feet | | Job # 86265,001 Site: CropSey |
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| r | Boring # B-5 MW# Street / of 2 |
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| | Project: AMA-Site TODESTIMATION |
| | Job # 85265,001 Site: Cr00, Seld |
| | Logged By: HRM Proj. Eng: DTS Edited By: |
| | Drilling Contractor: AES |
| | Drill Rig Type/Method: (100 probe |
| · · · · | Drillers Name: CHUCK / TOM |
| | Borehole Dlam./Drill Bit Type Total Depth 20 |
| | 1/2 ×5 Ref. Elev. 100 |
| | Hammer Wt: NA Drop: NA |
| | Start Time: 11/6 Date: 7/13/12 |
| HC=Hand Clear | Completion Time: // 45 Date: 7/13/12 |
| Sketch Map of Site Area With Drilling Locations | Backtill Time: Date: |
| Ci Ci Depth | 1st Water |
| PID/OVA Sampler Type & Depth Advance (in.) Advance (in.) Hecovered (In.) Hecovered (In.) Recovered siz | Casing Depth (ft.) |
| PID/OVA Sampler Type & Depth Blow Counts / (Advance (in.) Hecovered (in.) Recovered (in.) Advance filler Annulus Filler Sample Rec. Analyses Feet | Water Depth (fL) |
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| PID/OVA | Sampler Type | Blow Counts | Advance (In.) | Rcvd (In.) | | Casing | Annulus | Sample Rec. | Analysea | Feet | | Job # 8526,001 Site: Cropsler |
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| r | ; | Boring # B-6 MW# Street / at 3 |
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| | | Project: AMA |
| | | Job # 85265,001 Site: CropSly |
| | | Logged By: HRM Proj. Eng: D75 Edited By: |
| • | | Drilling Contractor: AES |
| | | Drill Rig Type/Method: 61600060 |
| | · | Drillers Name: Chuck/Tom |
| | | Borehole Diam./Drill Bit Type Total Depth |
| | | <u>//a X 5</u> Ref. Elev. |
| | | Hammer Wt: NA Drop: MA |
| he down along | | Start Time: 1/45 Date: 7/13/13 |
| HC=HandClear | | Completion Time: 12/5 Date: 7/3/12 |
| Sketch Map of Site Area With Drilling Lo | cations | Backtill Time: Date: |
| e In. | Depth | 1st Water Boring Depth (ft.) |
| PID/OVA Sampler Type & Depth Blow Counts / 6 In. Advance (In.) Recovered (In.) Casing Type & Siz | Jec | Casing Depth (ft.) |
| PID/OVA Sampler Type & Depth Blow Counts / Advance (in.) Recovered (in.) Recovered (in.) Recovered (in.) | Sample Rec. Analyses Feet | Water Depth (fL) |
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| | | | | | | | | | Project: AMA-Site Investigation |
| | | | | | | | | | Job # 85315,001 Site: (00560) |
| | | | | | | | | | Logged By: HPM Proj. Eng: OJS Edited By: |
| | | | | | | | | | Drilling Contractor: AES |
| | | | | | | | | | Drill Rig Type/Method: C1000000 |
| | | | | | | | | | Drillers Name: Chuck / Tom |
| | | | | | | | | | Borehole Diam./Drill Bit Type Total Depth 20' |
| | | | | | | | | | 1/2" x5" Ref. Elev. 100 |
| | | | | | | | | | Hammer Wt: NA Drop: NA |
| | | | | | | | | | Start Time: 1230 Date: 71311 |
| | HC = | Hord | Clean | ſ | | | | | Completion Time: 1400 Date: 71311 |
| | <u>s</u> | Sketch Map | of Site A | rea With Dr | rilling L | ocations | | | Backfill Time: Date: |
| | | <u>9</u> | | Size | | D | epth | | 1st Water Boring Depth (fL) |
| | Sampler Type & Depth | Blow Counts / 6 In. Advance (in.) | Recovered (in.) | Casing Type & | -iller | 90 | | | Casing Depth (ft.) |
| PID/OVA | npler Jepth | Blow Counts , Advance (in.) | DVer | L виј | Annulus Fliler | Sample Rec. Analyses | _ | | Water Depth (ft.) Time |
| PIC | | | | | | Sam Ana | Feel | | Date |
| 0.0 | <u>) HC</u> | 60' | O | D/A C | fice ficte | | | | |
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| <u>0,0</u> | m 5' | (ص" | 30" | | | <u> </u> | | | SP, Brawn, moist, med sand |
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| <u></u> | | - V 40" | 30" | | σŊ | | 10 - | | |
| <u>0.0</u> | 5' | <u> </u> | | | b | ₿-8 | | | |
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| A ler | | (in) | Ē | sul | e Rec. | 5 | | | Project: AMA |
| PID/OVA Sampler Type | Advance (In.) | Rcvd (In.) | Casing | Annulus | Sample Rec. | Analyses Fact | Lee | ſ | Job # 85265,001 Site: CropSQ4 |
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| | | | | | | | | | | | | Project: AMA-Site Investigation |
| | | | | | | | | | | | | Job # 85265,00/ Site: (100,000) |
| | | , | | | | | | | | | | Logged By: HRM Proj. Eng: DJS Edited By: |
| | | | | | | | | | | | | Drilling Contractor: $A \in S$ |
| | | | | | | | | | | | | Drill Rig Type/Method: Geoproble |
| | | | | | | | | | | • | | Drillers Name: Chuck / Torn |
| | | | | | | | | | | | | Borehole Dlam./Drill Bit Type Total Depth 20 |
| | | | | | | | | | | | | 1/2" × 5' Ref. Elev. 100 |
| | | | | | | | | | | | | Hammer Wt: NA Drop: NA |
| | | | | | | | | | | | | Start Time: 1200 Date: 7/13/12 |
| ц/ | <u>]</u> =H | $\sim r$ | | | / | | | | | | | Completion Time: //O) Date: 7/13/12 |
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| | 6d | Blow Counts / 6 in. | 1 | (in) | | θÅS | 9 | | | _ | | Boring Depth (ft.) |
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| PID/OVA | Sampler Type & Depth | low (| Advance (in.) | Recovered (In.) | | Casing Type & | Annulus Filler | Semolo Boo | Analyses | Feel | | Time 54, |
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| | 5' | | | | | | | | , | | · | SP. Brawn, moist, med sand No adors, wa staining |
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| | -{ | | $\left \right $ | | - | S | + | IK | -4 | 8 — | ╞ | med sond / some gravel |
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| r | Bóring # MW# 3 Sheet I of O |
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| | Project: AMA-Site Investigation |
| | Job # 852(05,001 Site: Cra)Sel |
| | Logged By: I-\ RM Proj. Eng: DTS Edited By: |
| | Drilling Contractor: AES |
| | Drill Rig Type/Method: GROODDR |
| | Drillers Name: CINCLCK /TOM |
| | Borehole Dlam./Drill Bit Type Total Depth 20 |
| | 112 × 5 |
| | Hammer Wt: NA Drop: NA |
| | Start Time: 1000 Date: 7112/12 |
| HC=Hand Clear | Completion Time: 1130 Date: 71212 |
| Sketch Map of Site Area With Drilling Locations | Backtill Time: Date: |
| C Depth | 1st Water |
| | Boring Depth (ft.) |
| 2VA DVA Court 1 D D D D D D D D D D D D D D D D D D | Water Depth (fL) |
| PID/OVA Sampler Type & Depth Blow Counts / 6 In.) Advance (In.) Recovered (In.) Recovered (In.) Recovered (In.) Recovered (In.) Recovered (In.) Recovered (In.) | Time Ni. |
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| PID/OVA | ampler | Type | Blow Counts | | Advance (In.) | Rcvđ (in.) | | Casing | Annulus | Sample Rec. | Analyse | Feet | | Proje | | <u> -</u> - - | <u>1</u> 5.(| <u></u> ງ _ກ ູ | Site | $\overline{0}$ | 5 | 2 | ee | • |
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| | | | Project: Cropsey | Boring ID MW-1S | | |
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| | | | Brooklyn, NY | Sheet 1 of 1 | | |
| Boring Location: Fenced are | | dry cleaner | Project Manager: R. Costanzo | Project Number: 85265 | | |
| Ground Elevation: | 22.9 | | Dated Started: 7/13/11 | Date completed: 7/13/11 | | |
| | | | Drill Type: Direct Push for soil cores; Hollow | | | |
| Depth to First Water: | 18 | | stem auger for well installation | Borehole Dia: 4.25" | | |
| Depth to Static Water: | 19.7 | | Drill Rig Type: | Geoprobe | | |
| Top of Casing Elevation: | 22.6 | | Drilling Company: | Zebra | | |
| 1 8 | Samp | ler | Driller's Name: | Chuck | | |
| Size: 2" I.D. PVC | Туре: | <u></u> | APEX Representative: | HRM | | |
| Screen: 0.010-inch Slot PVC | Lammar: | NA | Owner/Client Rep.: | AMA Cropsey | | |
| Sand: #2 Morie | Fall: | NA | Owner/Chent Rep | AMA Clopsey | | |
| | | NA | | | | |
| ti Control No. | PID (ppm) USCS Class. | D | escription of Soils | Well Construction | | |
| 0 | 0 | Hand Clear | | Locking Cap at grade | | |
| 5 2.5 10 2.5 MW-1-10'-15' 15 4 MW-1-15'-20' 20 25 30 | 0 SP 0 0 0 | Brown medium grit sand, moist, no od Groundwater encountered at 18' bgs | ors, no staining | Slotted screened interval from 18-28' bgs | | |
| Locking Cap | | Morie Sand #2 | | | | |
| Screened interval | | Bentonite seal | | | | |
| | | Grout | | | | |

| | | | | | | Project: Cropsey | Boring ID MW | /-1I |
|-----------------|-------------------|-------------------------------|------------|----------------|---|--|---------------------------------------|---------|
| Danin | - T 4 | Service Constant | | £ | f 1 | Brooklyn, NY | Sheet 1 of 1 Project Number: 85265 | |
| | g Local d Elev | | 22.6 | | leaner fenced area | Project Manager: R. Costanzo Dated Started: 1/19/2016 | Date completed: 1/19/1 | |
| Groun | a Elev | ation: | 22.0 | | | Drill Type: Direct Push for soil cores; Hollow | Date completed: 1/19/1 | 0 |
| Depth | to Firs | st Water: | | | | stem auger for well installation | Borehole Dia: 4.25" | |
| | | tic Water: | 19.2 | | | Drill Rig Type: | Geoprobe | |
| | | g Elevation: | 22.2 | | | Drilling Company: | Zebra | |
| | | 0 | | Samp | ler | Driller's Name: | Evan | |
| | 2" I.D. | | Type: | | Encore | APEX Representative: | Kristen Main/Daniel Ko | opec |
| | | 0-inch Slot PVC | | ner: | NA | Owner/Client Rep.: | AMA Cropsey | |
| Sand: | #2 Mo | rie | Fall: | 1 | NA | | | |
| ч — | ery | | PID (ppm) | s .: | | | | |
| Depth (feet) | Recovery | Sample I.D. or Control No. | Id) | USCS Class. | D | escription of Soils | Well Constr | ruction |
| D O | Rec | Control No. | DIG | рo | | | | |
| 0 | N/A | N/A | 0 | SM | 1':Topsoil/grass, black, some sand, dry | 1 | Locking Cap at grade | |
| E | | | | | 2-3':Brown silty clay, little to no sand, | | 010 | |
| L . | | | | | | | | |
| ┝ | | | | | 3-5':Brown medium-coarse sand, 10-2 | 0% gravel, moist | | |
| - 5 | 3 | N/A | * | | | | | |
| Ľ | | | | | | | | |
| E . | | | | SW | 7': Brown medium-coarse sand, little to | o no fines, dry | | |
| F | | | | | | 1 | | |
| - 10 | 2.7 | N/A | * | | 9': Brown fine-medium sand, dry, som | e graveis | | |
| - 10 | 2.7 | 1011 | | SW | 11': Brown medium to coarse sand, litt | tle to no fines, 5-10% gravels, dry | | |
| E | | | | SW | 12': Brown medium to coarse sand, litt | tle to no fines | | |
| F | | | | | 13': SAA, 0-5% gravels | | | |
| - 15 | 4 | | * | | | | | |
| - 13 | 4 | | | SW | 16': Brown fine to medium sand, 5-10 | % gravels | | |
| F | | | | SW | | nd, slightly moist, 0-5% gravels, 5-10% fines | | |
| E | | | | | | | | |
| | | | <u>ب</u> | | | | | |
| 20 | 4 | MW-11-20 | * | | | | | |
| - | | | | SW | 22': Brown medium to coarse sand, 0-: | 5% gravel, wet | | |
| L | | | | | Groundwater encountered at ~23' bgs | - | | |
| | | | | | | | | |
| 25 | 5 | MW-11-25 | * | SW | Brown medium to coarse sand, 10-15% | % gravels, 5-10% fines, wet | | |
| F | | | | | | | | |
| Ľ | | | | | | | | |
| L | _ | | | | | | | |
| 30 | 5 | MW-1I-30 | * | SW | Brown medium to coarse sand, 0-5% f | ines, 0-5% gravels. Wet | | |
| F | | | | | | | | |
| F | | | | | | | | |
| _ | | | | SP | 34-35': Brown fine sand | | | |
| 35 | 5 | MW-1I-35 | * | SP | Brown/light red medium to coarse san | d, wet | | |
| F | | | | | | | | |
| Ľ | | | | | | | | |
| L . | | | | | | | | |
| 40 | 5 | MW-1I-40 | * | SP | SAA | | | |
| F | | | | | | | Slotted screened interval | |
| Ē | | | | | | | from 40-50' bgs | |
| Ľ | | | | | | | | |
| 45 | 5 | MW-1I-45 | * | SP | SAA | | | |
| F | | | | | | | | |
| Ľ | | | | SM | 46.5': Brown medium sand, 10-15% fi | nes, wet | | |
| F | | | | a | | | | |
| 50 | | MW-1I-50 | * | SM | SAA | | | |
| | | I | L | 1 | I | | I | |
| | Locking | g Cap | | | Morie Sand #2 | | | |
| | Saraar | d interval | | 1 | Bentonite seal | | | |
| | Screene | a micivai | | | Demolific scal | | | |
| | | | | | Grout | | | |
| 1 | * | Outside temp. duri | ng drilliı | ng was to | oo cold for the PID to work properly | | | |

| | | | Project: Cropsey | Boring ID MW-2 |
|---|-----------------------------|---|--|--|
| | | | Brooklyn, NY | Sheet 1 of 1 |
| Boring Location: Courtyard | behind stripn | nall | Project Manager: R. Costanzo | Project Number: 85265 |
| Ground Elevation: | 22.7 | | Dated Started: 7/13/12 | Date completed: 7/13/12 |
| | | | Drill Type: Direct Push for soil cores; Hollow | 1 |
| Depth to First Water: | 18 | | stem auger for well installation | Borehole Dia: 4.25" |
| Depth to Static Water: | 19.3 | | Drill Rig Type: | Geoprobe |
| Top of Casing Elevation: | 22.3 | | Drilling Company: | Zebra |
| Top of Casing Elevation. | Samp | 1 | Driller's Name: | Chuck |
| | | ler | | HRM |
| Size: 2" I.D. PVC Screen: 0.010-inch Slot PVC | Туре: | NT 4 | APEX Representative: | |
| | | NA | Owner/Client Rep.: | AMA Cropsey |
| Sand: #2 Morie | Fall: | NA | | |
| (1) (1) (1) (1) (1) (1) (1) (1) | PID (ppm) USCS Class. | D | escription of Soils | Well Construction |
| 0 0 | 0 | Hand Clear | | Locking Cap at grade |
| 5 4 10 3.5 MW-2-10'-15' 15 4.5 MW-2-15'-20' 20 25 | 0 SP 0 | Brown medium sand, moist, no odors, Groundwater encountered at 18' bgs | no staining | Slotted screened interval from 18-28' bgs |
| Locking Cap | | Morie Sand #2 | | |
| Screened interval | | Bentonite seal | | |
| | | Grout | | |

| | | | Project: Cropsey | Boring ID MW-3 |
|--|-----------------------------|---|--|--|
| | | | Brooklyn, NY | Sheet 1 of 1 |
| Boring Location: Courtyard | | nall | Project Manager: R. Costanzo | Project Number: 85265 |
| Ground Elevation: | 23.1 | | Dated Started: 7/13/12 | Date completed: 7/13/12 |
| | | | Drill Type: Direct Push for soil cores; Hollow | |
| Depth to First Water: | 18 | | stem auger for well installation | Borehole Dia: 4.25" |
| Depth to Static Water: | 19.7 | | Drill Rig Type: | Geoprobe |
| Top of Casing Elevation: | 22.6 | | Drilling Company: | Zebra |
| | Samp | ler | Driller's Name: | Chuck |
| Size: 2" I.D. PVC | Туре: | | APEX Representative: | HRM |
| Screen: 0.010-inch Slot PVC | Hammer: | NA | Owner/Client Rep.: | AMA Cropsey |
| Sand: #2 Morie | Fall: | NA | * | * * |
| Line for the second sec | PID (ppm) USCS Class. | D | escription of Soils | Well Construction |
| 0 0 | 0 | Hand Clear | | Locking Cap at grade |
| 5 4 10 4.5 MW-3-10'-15' 15 4.5 MW-3-15'-20' 20 25 30 | 0 SP 0 | Brown medium sand, moist, no odors, Groundwater encountered at 18' bgs | no staining | Slotted screened interval from 18-28' bgs |
| Locking Cap | | Morie Sand #2 | | |
| Screened interval | | Bentonite seal | | |
| | | Grout | | |

| | | | | | | During to Comment | Boring ID MW | 7 48 | | |
|-----------------|----------------|-------------------------------|------------|----------------|--|--|-------------------------------------|---------|--|--|
| | | | | | | Project: Cropsey | Boring ID MW Sheet 1 of 1 | /-48 | | |
| Domina | Looot | ion. Sidowall | on the | | of Cropsey Ave and 20th Ave | Brooklyn, NY Project Manager: R. Costanzo | Project Number: 85265 | | | |
| Ground | | | <u>22</u> | corner | of Cropsey Ave and 20th Ave | Dated Started: 1/19/2016 | | | | |
| Groun | a Eleva | ation: | 22 | | | Drill Type: Direct Push for soil cores; Hollow | Date completed: 1/19/1 | 0 | | |
| Donth | to Eina | + Watan | 23 | | | | Darahala Dia, 4.25" | | | |
| | | t Water: | | | | stem auger for well installation | Borehole Dia: 4.25" | | | |
| | | ic Water: | 18.6 | | | Drill Rig Type: | Geoprobe Zebra | | | |
| 100 01 | Casin | g Elevation: | 23.1 | Sampl | la n | Drilling Company: Driller's Name: | Evan | | | |
| Size: 2 | " I D | DVC | | Sampi | Encore | APEX Representative: | Evan Kristen Main/Ashlyn Norberg | | | |
| | | -inch Slot PVC | Type: | or | NA | Owner/Client Rep.: | AMA Cropsey | orberg | | |
| Sand: | | | Fall: | ICI. | NA | Owner/Chent Rep | AWA Cropsey | | | |
| Sand. | | | | | | | | | | |
| Depth (feet) | Recovery | Sample I.D. or Control No. | PID (ppm) | USCS Class. | D | Description of Soils | Well Constr | ruction | | |
| D) | Re | Control No. | PID | 0 | | | | | | |
| 0 | N/A | N/A | 0 | | 0-0.5': Concrete | | Locking Cap at grade | | | |
| | | | | | 0.5-5': Reddish brown medium to coar | se sand, dry | 6 | | | |
| [| | | | | | | | | | |
| L | | | | | | | | | | |
| - 5 | N/A | N/A | N/A | | | | | | | |
| - 3 | IN/A | 18/24 | IN/A | | | | | | | |
| - | | | | | | | | | | |
| _ | | | | | | | | | | |
| | | | | | | | | | | |
| 10 | 10 N/A N/A N/A | | | | | | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| 15 | N/A | N/A | N/A | | | | | | | |
| _ | | | | | | | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| 20 | 2 | | * | SP | Reddish brown medium to coarse sand | l. wet | | | | |
| | - | | | | second of the mount to course said | | Slotted screened interval | | | |
| _ | | | | | | | from 17-27' bgs | | | |
| _ | | MW-4S-23.5 | * | | Groundwater encountered at ~23 ft | | | | | |
| - | | | | _ | | | | | | |
| 25 | 3.5 | | * | SP | SAA | | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| 30 | | MW-4S-30 | * | SW | Brown medium-coarse sand, trace grav | vel, trace mica, wet | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| - 35 | 3.5 | N/A | * | | | | | | | |
| | 2.00 | | | SW | 36.5-37.5': SAA | | | | | |
| | | | | SM | 37.5-38': Fine-medium grit browm san | ıd | | | | |
| _ | | | | SP | 38-40': Medium to coarse sand | | | | | |
| | | 101/10/10 | * | | | | | | | |
| 40 | | MW-4S-40 | * | | | | | | | |
| | L | 1 | | | <u> </u> | | 1 | | | |
| | Locking | g Cap | | | Morie Sand #2 | | | | | |
| | Screene | d interval | | | Bentonite seal | | | | | |
| | | | | | | | | | | |
| 1 | * | Outside temp. duri | ng drillin | g was to | Grout Grout for the PID to work properly | | | | | |
| | | 1 | | - | 1 1 | | | | | |

| Brocklyn, NV Short I of 1 Brong Location: Sidewalk on the corner of Cropsey Ave and 20th Ave Ground Elevation: Project Manager: R. Costaro? Duded Started: Duded Started: Dude Star | | | | | | | Project: Cropsey | Boring ID MV | V-4I | | |
|---|---------|---------|--------------------|------------|-----------|---------------------------------------|--|---------------------------|---------|--|--|
| Group Le varion: 2 Datel Started: pate ompleted: Dopth to First Water: 16 Datel Started: Date ompleted: To of Costang Elevation: 21.6 Defiling: Company: Zehra Starte: Defiling: Company: Zehra Concolo Dia: 42.5" Starte: Datel Starte: Defiling: Company: Zehra Starte: Differ: Starte: Starte: Starte: Starte: Differ: Name Version Starte: Starte: Name Na Na Na Starte: Control No. Starte: Na Na Starte: Control No. Starte: Na Na Starte: Control No. Starte: Starte: Na Starte: Control No. Starte: Starte: Na Na Na Na Na Na Na Starte: Control No. Starte: Starte: Na Starte: Na Na | | | | | | | Brooklyn, NY | Sheet 1 of 1 | | | |
| Group Le varion: 2 Datel Started: pate ompleted: Dopth to First Water: 16 Datel Started: Date ompleted: To of Costang Elevation: 21.6 Defiling: Company: Zehra Starte: Defiling: Company: Zehra Concolo Dia: 42.5" Starte: Datel Starte: Defiling: Company: Zehra Starte: Differ: Starte: Starte: Starte: Starte: Differ: Name Version Starte: Starte: Name Na Na Na Starte: Control No. Starte: Na Na Starte: Control No. Starte: Na Na Starte: Control No. Starte: Starte: Na Starte: Control No. Starte: Starte: Na Na Na Na Na Na Na Starte: Control No. Starte: Starte: Na Starte: Na Na | Boring | Locat | ion: Sidewalk | on the o | corner | of Cropsey Ave and 20th Ave | | Project Number: 85265 | 5 | | |
| Depth to First Water: Image for weal installation Borchole Dia: 4.25" Depth to Strict Water: 18.6 Drill Ng Company: Zebra To of Casing Livation: 21.6 Drill Ng Company: Zebra Streep: D107 info Stot PVC Type: First Water: NA Streep: D107 info Stot PVC Type: First Water: NA Streep: D107 info Stot PVC Type: First Water: NA Streep: D107 info Stot PVC Type: NA Owner/Client Rep:: AMA Cropsey Streep: Control No. Editary Streep: NA Owner/Client Rep:: AMA Cropsey 9 NA NA Info Stot PVCHarme: NA Description of Soils Well Construction 10 NA NA NA Info Stot PVCHarme: Stote Output Na Stote Output Na Stote Output Na 10 NA NA NA Stote Output Na Stote Output Na Stote Output Na Stote Output Na 10 NA NA NA Stote Output | Ground | d Eleva | ation: | | | 1 2 | | | | | |
| Depth to First Water: Istem auger for well installation Boocholo Dia: 4.28" Depth to Statut Water: 1.6 Diali Rig Type:: Geoprole Size: 2"10. PVC Type: Incore APER Keynsentative: Steve Cotoms/Ashlyn Norberg: Size: 2"10. PVC Type: Incore APER Keynsentative: Steve Cotoms/Ashlyn Norberg: State: 2"10. PVC Type: Incore APER Keynsentative: Steve Cotoms/Ashlyn Norberg: State: 2"10. PVC Type: NA Owner/Client Rep:: APAC ropsey State: 2"Morie Fall: NA Owner/Client Rep:: APAC ropsey State: 4"2 Morie Fall: NA Owner/Client Rep:: APAC ropsey State: 4"2 Morie Fall: NA Owner/Client Rep:: APAC ropsey State: 4"2 Morie Fall: NA Owner reclinet Rep:: APAC ropsey State: 4"2 Morie Fall: NA Owner roccurse state dy 10 NA NA NA Index of the proven medium to coarse state, wet State determed internof 20 2 | | | | | | | Drill Type: Direct Push for soil cores: Hollow | | | | |
| Depth to Sudi: Water: 18.6 Delt Rig Type: Geoprobe Top of Casing Elevation: 21.6 Drilling Company: Zeba Server: DIU-inch Slot PVC Tamuer: NA Server: Server: MAA Cropsey Server: DIU-inch Slot PVC Tamuer: NA Owner/Client Rep: AMA Cropsey Sand: #22 Morie NA Owner/Client Rep: AMA Cropsey Server: Server: NA 9 \$\frac{1}{28}\$ Sample LD. or Control No. \$\frac{1}{28}\$ \$\frac{2}{29}\$ \$\frac{1}{28}\$ Server: Locking Cip at grade 10 NA NA 0 0.5.5?. Redifish hown medium to coarse said, dry Locking Cip at grade Locking Cip at grade 10 NA NA NA \$\frac{1}{28}\$ Server: Server: <td< td=""><td>Donth</td><td>to Eina</td><td>+ Watan</td><td></td><td></td><td></td><td colspan="3"></td></td<> | Donth | to Eina | + Watan | | | | | | | | |
| Top of Cashy Elevation: 21.6 Defining Company: Zohna Zohna Size: 27 10. PVC Type: Facore APEX Representative: Size: Control Aving | Depin | to Firs | t water: | 10.6 | | | | | | | |
| Size: 2* ID. PVC Type: Description Evan Screet: 0100-inch Stor PVC (Hammer: NA AA Owner/Client Rep: AMA Cropsey Sind: 22 Marie Full: NA Owner/Client Rep: AMA Cropsey Sind: 22 Marie Full: NA Owner/Client Rep: AMA Cropsey Sind: 22 Marie Full: NA Owner/Client Rep: AMA Cropsey Sind: 22 Marie Full: NA Owner/Client Rep: AMA Cropsey Sind: 22 Marie Full: NA Owner/Client Rep: AMA Cropsey Sind: 22 Marie Full: NA Owner/Client Rep: AMA Cropsey Sind: 24 Marie Full: NA Owner/Client Rep: Control No. Sind: 42 Marie Full: NA Description of Soils Well Construction Sind: 42 Marie Full: NA NA Soil: Control No. Sind: 5 N/A NA NA NA Soil: Control No. Soil: Sind: 5 N/A NA NA NA Soil: Control No. Soil: Sind: 5 N/A NA NA NA Soil: Control No. Soil: Sind: 5 N/A NA NA Soil: Groundwater encountered at -23 ft </td <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td colspan="4"></td> | | | | | | | | | | | |
| Size: 2:10. PVC Type: Encore APEX Representative: Steve Corron-Aships Norberg Sterie: 0.01-inde blor PC Hammer: NA OmerClient Rep.: AMA Cropsey Sterie: 0.01-inde blor PC Hammer: NA OmerClient Rep.: AMA Cropsey 9 9 Sample I.D. or Control No. 9 9 9 Sample I.D. or Control No. 9 9 9 NA Well Construction 1 NA NA 0 0.55': Sealiab brown medium to coarse sand, dry Lacking Cap at grade 0.55': Sealiab brown medium to coarse sand, dry Lacking Cap at grade 0.55': Sealiab brown medium to coarse sand, wet 0.55': Sealiab brown medium to coarse sand, wet 0.55': Sealiab brown medium to coarse sand, wet Stere Sealiab brown medium to coarse sand, wet Stere Sealiab brown medium to coarse sand, wet Stere Sealiab brown medium coarse sand, wet Stere Sealiab brown medium to coarse sand, wet Stere Sealiab brown sand Ste | Top of | 'Casin | g Elevation: | 21.6 | | | | Zebra | | | |
| Sieven: OII-vinck Sion PVC Harmer: NA Ovner/Client Rep.: AMA Cropsey 6 1 V Avian Via | | | | | Sampl | ler | Driller's Name: | | | | |
| Sieven: OII-vinck Sion PVC Harmer: NA Ovner/Client Rep.: AMA Cropsey 6 1 V Avian Via | Size: 2 | " I.D. | PVC | Type: | | Encore | APEX Representative: | Steve Cotrone/Ashlyn | Norberg | | |
| Sand: #2 Morie Full: NA If g g g g g g g g g Sample LD, or g g g g g g g g g g g g g g g g g g | | | | Hamm | ner: | | | | U | | |
| up by Sample LD. or Control No. Sty Sty Description of Soils Well Construction 0 N/A N/A 0 0 0.9.9.1 Concrete 0.5.9.1 Concrete Locking Cap at grade 5 N/A N/A N/A N/A N/A 0 0.5.9.1 Concrete Locking Cap at grade 10 N/A N/A N/A N/A N/A Locking Cap at grade Locking Cap at grade 20 2 N/A N/A N/A N/A Readdab brown medium to coarse sand, dry Locking Cap at grade 20 2 + + SP Readdab brown medium to coarse sand, wet Crossnowneed at23 ft 20 2 3.5 N/A N/A SP Readdab brown medium to coarse sand, wet 30 MW-45-20.5 + SP SAA 30 MW-45-30 + SP SAA 31 3.5 N/A + SP 32 3.5 N/A + SP 33 3.5 N/A + SP 34 3.5 N/A + SP 35 X + SP SAA 36 3.5 N/A | | | | | | | - · · · · · · · · · · · · · · · · · · · | | | | |
| 0 N/A N/A 0 0.45* Concrete 0.5*5* Redish brown medium to coarse sand, dry 1 N/A N/A N/A N/A 10 N/A N/A N/A 115 N/A N/A N/A 120 2 * SP 20 2 * SP 20 2 * SP 20 2 * SP 21 XW-45-23.5 * Groundwater encountered at -23 ft 22 3.5 N/A SW Brown medium-coarse sand, trace gravel, trace mica, wet 30 MW-45-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 N/A * SW 35-37.5: SAA 40 MW-45-40 * SW 35-37.5: SAA 40 MW-45-40 * Morie Sund #2 < | Suna. I | | | | 1 | 1111 | | | | | |
| 0 N/A N/A 0 0.45* Concrete 0.5*5* Redish brown medium to coarse sand, dry 1 N/A N/A N/A N/A 10 N/A N/A N/A 115 N/A N/A N/A 120 2 * SP 20 2 * SP 20 2 * SP 20 2 * SP 21 XW-45-23.5 * Groundwater encountered at -23 ft 22 3.5 N/A SW Brown medium-coarse sand, trace gravel, trace mica, wet 30 MW-45-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 N/A * SW 35-37.5: SAA 40 MW-45-40 * SW 35-37.5: SAA 40 MW-45-40 * Morie Sund #2 < | 4 0 | ery | G 1 1 D | m | <i>s</i> | | | | | | |
| 0 N/A N/A 0 0.45* Concrete 0.5*5* Redish brown medium to coarse sand, dry 1 N/A N/A N/A N/A 10 N/A N/A N/A 115 N/A N/A N/A 120 2 * SP 20 2 * SP 20 2 * SP 20 2 * SP 21 XW-45-23.5 * Groundwater encountered at -23 ft 22 3.5 N/A SW Brown medium-coarse sand, trace gravel, trace mica, wet 30 MW-45-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 N/A * SW 35-37.5: SAA 40 MW-45-40 * SW 35-37.5: SAA 40 MW-45-40 * Morie Sund #2 < | eet | Ň | | Id) | ase | D | Description of Soils | Well Const | ruction | | |
| 0 N/A N/A 0 0.45* Concrete 0.5*5* Redish brown medium to coarse sand, dry 1 N/A N/A N/A N/A 10 N/A N/A N/A 115 N/A N/A N/A 120 2 * SP 20 2 * SP 20 2 * SP 20 2 * SP 21 XW-45-23.5 * Groundwater encountered at -23 ft 22 3.5 N/A SW Brown medium-coarse sand, trace gravel, trace mica, wet 30 MW-45-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 N/A * SW 35-37.5: SAA 40 MW-45-40 * SW 35-37.5: SAA 40 MW-45-40 * Morie Sund #2 < | Ű D | ec | Control No. | Ω | 5 D | | 1 | | | | |
| 1 NA NA NA NA NA NA 10 NA NA NA NA Image: Constraint of the constraint | | | | IJ | | | | | | | |
| s NA | 0 | N/A | N/A | 0 | | | | Locking Cap at grade | | | |
| 10 N/A N/A N/A I 15 N/A N/A N/A I 16 N/A N/A N/A I 17 N/A N/A N/A I 18 N/A N/A N/A I 19 N/A N/A N/A I 19 N/A N/A SP Reddish brown medium to course sand, wet 10 MW-48-23.5 * Groundwater encountered at -23 ft Groundwater encountered at -23 ft 10 N/A N/A SP SA SA 10 MW-48-30 * SP SA 10 MW-48-30 * SP SA- 10 MW-48-40 * SP SA- 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * * SP 10 MW-48-40 * * SP 10 MW-48-40 * * < | | | | | | 0.5-5': Reddish brown medium to coar | rse sand, dry | | | | |
| 10 N/A N/A N/A I 15 N/A N/A N/A I 16 N/A N/A N/A I 17 N/A N/A N/A I 18 N/A N/A N/A I 19 N/A N/A N/A I 19 N/A N/A SP Reddish brown medium to course sand, wet 10 MW-48-23.5 * Groundwater encountered at -23 ft Groundwater encountered at -23 ft 10 N/A N/A SP SA SA 10 MW-48-30 * SP SA 10 MW-48-30 * SP SA- 10 MW-48-40 * SP SA- 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * * SP 10 MW-48-40 * * SP 10 MW-48-40 * * < | | | | | | | | | | | |
| 10 N/A N/A N/A I 15 N/A N/A N/A I 16 N/A N/A N/A I 17 N/A N/A N/A I 18 N/A N/A N/A I 19 N/A N/A N/A I 19 N/A N/A SP Reddish brown medium to course sand, wet 10 MW-48-23.5 * Groundwater encountered at -23 ft Groundwater encountered at -23 ft 10 N/A N/A SP SA SA 10 MW-48-30 * SP SA 10 MW-48-30 * SP SA- 10 MW-48-40 * SP SA- 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * * SP 10 MW-48-40 * * SP 10 MW-48-40 * * < | | | | | | | | | | | |
| 10 N/A N/A N/A I 15 N/A N/A N/A I 16 N/A N/A N/A I 17 N/A N/A N/A I 18 N/A N/A N/A I 19 N/A N/A N/A I 19 N/A N/A SP Reddish brown medium to course sand, wet 10 MW-48-23.5 * Groundwater encountered at -23 ft Groundwater encountered at -23 ft 10 N/A N/A SP SA SA 10 MW-48-30 * SP SA 10 MW-48-30 * SP SA- 10 MW-48-40 * SP SA- 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * SP SA-37.5: SA A 10 MW-48-40 * * SP 10 MW-48-40 * * SP 10 MW-48-40 * * < | | | | | | | | | | | |
| 15 N/A N/A N/A N/A 20 2 * SP Reddish brown medium to coarse sand, wet 20 2 * SP Reddish brown medium to coarse sand, wet 25 3.5 * SP Groundwater encountered at ~23 ft 26 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 32 3.5 N/A * SW 36.5-37.5': SAA 33 4.0': MW-4S-40 * * SW 36.5-37.5': SAA 34 3.5 N/A * SW 36.5-37.5': SAA 35 N/A * SW 36.5-37.5': SAA SM 36 MW-4S-40 * Wrie Sand #2 Mrie Sand #2 | 5 | N/A | N/A | N/A | | | | | | | |
| 15 N/A N/A N/A N/A 20 2 * SP Reddish brown medium to coarse sand, wet 20 2 * SP Reddish brown medium to coarse sand, wet 25 3.5 * SP Groundwater encountered at ~23 ft 26 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 32 3.5 N/A * SW 36.5-37.5': SAA 33 4.0': MW-4S-40 * * SW 36.5-37.5': SAA 34 3.5 N/A * SW 36.5-37.5': SAA 35 N/A * SW 36.5-37.5': SAA SM 36 MW-4S-40 * Wrie Sand #2 Mrie Sand #2 | | | | | | | | | | | |
| 15 N/A N/A N/A N/A 20 2 * SP Reddish brown medium to coarse sand, wet 20 2 * SP Reddish brown medium to coarse sand, wet 25 3.5 * SP Groundwater encountered at ~23 ft 26 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 32 3.5 N/A * SW 36.5-37.5': SAA 33 4.0': MW-4S-40 * * SW 36.5-37.5': SAA 34 3.5 N/A * SW 36.5-37.5': SAA 35 N/A * SW 36.5-37.5': SAA SM 36 MW-4S-40 * Wrie Sand #2 Mrie Sand #2 | | | | | | | | | | | |
| 15 N/A N/A N/A N/A 20 2 * SP Reddish brown medium to coarse sand, wet 20 2 * SP Reddish brown medium to coarse sand, wet 25 3.5 * SP Groundwater encountered at ~23 ft 26 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 32 3.5 N/A * SW 36.5-37.5': SAA 33 4.0': MW-4S-40 * * SW 36.5-37.5': SAA 34 3.5 N/A * SW 36.5-37.5': SAA 35 N/A * SW 36.5-37.5': SAA SM 36 MW-4S-40 * Wrie Sand #2 Mrie Sand #2 | | | | | | | | | | | |
| 15 N/A N/A N/A N/A 20 2 * SP Reddish brown medium to coarse sand, wet 20 2 * SP Reddish brown medium to coarse sand, wet 25 3.5 * SP Groundwater encountered at ~23 ft 26 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 31 3.5 N/A * SW 36.5-37.5': SAA 32 3.5 N/A * SW 36.5-37.5': SAA 33 4.0': MW-4S-40 * * SW 36.5-37.5': SAA 34 3.5 N/A * SW 36.5-37.5': SAA 35 N/A * SW 36.5-37.5': SAA SM 36 MW-4S-40 * Wrie Sand #2 Mrie Sand #2 | | | | | | | | | | | |
| 20 2 ** SP Reddish brown medium to coarse sand, wet 25 3.5 MW-4S-23.5 ** SP Solution and the coarse sand, wet 30 . MW-4S-30 ** SP SAA 30 . MW-4S-30 ** SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 31 3.5 NA * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 40 . MW-4S-40 * . SW Brown medium to coarse sand 40 . MW-4S-40 * 50 40 50 60 | 10 | N/A | N/A | N/A | | | | | | | |
| 20 2 ** SP Reddish brown medium to coarse sand, wet 25 3.5 MW-4S-23.5 ** SP Solution and the coarse sand, wet 30 . MW-4S-30 ** SP SAA 30 . MW-4S-30 ** SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 31 3.5 NA * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 40 . MW-4S-40 * . SW Brown medium to coarse sand 40 . MW-4S-40 * 50 40 50 60 | | | | | | | | | | | |
| 20 2 ** SP Reddish brown medium to coarse sand, wet 25 3.5 MW-4S-23.5 ** SP Solution and the coarse sand, wet 30 . MW-4S-30 ** SP SAA 30 . MW-4S-30 ** SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 31 3.5 NA * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 40 . MW-4S-40 * . SW Brown medium to coarse sand 40 . MW-4S-40 * 50 40 50 60 | | | | | | | | | | | |
| 20 2 ** SP Reddish brown medium to coarse sand, wet 25 3.5 MW-4S-23.5 ** SP Solution and the coarse sand, wet 30 . MW-4S-30 ** SP SAA 30 . MW-4S-30 ** SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 31 3.5 NA * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 40 . MW-4S-40 * . SW Brown medium to coarse sand 40 . MW-4S-40 * 50 40 50 60 | | | | | | | | | | | |
| 20 2 | | | | | | | | | | | |
| 25 3.5 MW-4S-23.5 * SP Groundwater encountered at ~23 ft 30 A SP SAA 30 MW-4S-30 * SP 31 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 32 3.5 NA SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 NA * SW 34 MW-4S-40 * SW 36-37.5: SAA 35 NA * SW 36-37.5: SAA 36 MW-4S-40 * SW 36-37.5: SAA 37.5-383: Fine-medium grit brown sand SP 38-40: Medium to coarse sand 38 MW-4S-40 * SW 36-37.5: SAA 39 Beine-medium grit brown sand SP 38-40: Medium to coarse sand 39 Verie Sand #2 Merie Sand #2 Merie Sand #2 | 15 | N/A | N/A | N/A | | | | | | | |
| 25 3.5 MW-4S-23.5 * SP Groundwater encountered at ~23 ft 30 A SP SAA 30 MW-4S-30 * SP 31 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 32 3.5 NA SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 NA * SW 34 MW-4S-40 * SW 36-37.5: SAA 35 NA * SW 36-37.5: SAA 36 MW-4S-40 * SW 36-37.5: SAA 37.5-383: Fine-medium grit brown sand SP 38-40: Medium to coarse sand 38 MW-4S-40 * SW 36-37.5: SAA 39 Beine-medium grit brown sand SP 38-40: Medium to coarse sand 39 Verie Sand #2 Merie Sand #2 Merie Sand #2 | | | | | | | | | | | |
| 25 3.5 MW-4S-23.5 * SP Groundwater encountered at ~23 ft 30 A SP SAA 30 MW-4S-30 * SP 31 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 32 3.5 NA SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 NA * SW 34 MW-4S-40 * SW 36-37.5: SAA 35 NA * SW 36-37.5: SAA 36 MW-4S-40 * SW 36-37.5: SAA 37.5-383: Fine-medium grit brown sand SP 38-40: Medium to coarse sand 38 MW-4S-40 * SW 36-37.5: SAA 39 Beine-medium grit brown sand SP 38-40: Medium to coarse sand 39 Verie Sand #2 Merie Sand #2 Merie Sand #2 | | | | | | | | | | | |
| 25 3.5 MW-4S-23.5 * SP Groundwater encountered at ~23 ft 30 A SP SAA 30 MW-4S-30 * SP 31 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 32 3.5 NA SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 NA * SW 34 0.5 NA SW 36-37.5: SAA 35 NA * SW 36-37.5: SAA 36 MW-4S-40 * SW 36-37.5: SAA 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 36 MW-4S-40 * SW 36-37.5': SAA 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 38 Wine Sand #2 Mine Sand #2 Mine Sand #2 | | | | | | | | | | | |
| 25 3.5 MW-4S-23.5 * SP Groundwater encountered at ~23 ft 30 A SP SAA 30 MW-4S-30 * SP 31 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 32 3.5 NA SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 NA * SW 34 0.5 NA SW 36-37.5: SAA 35 NA * SW 36-37.5: SAA 36 MW-4S-40 * SW 36-37.5: SAA 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 36 MW-4S-40 * SW 36-37.5': SAA 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 38 Wine Sand #2 Mine Sand #2 Mine Sand #2 | | | | | | | | | | | |
| 25 3.5 MW-4S-23.5 * SP Groundwater encountered at ~23 ft 30 A SP SAA 30 MW-4S-30 * SP 31 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet 32 3.5 NA SW Brown medium-coarse sand, trace gravel, trace mica, wet 33 3.5 NA * SW 34 0.5 NA SW 36-37.5: SAA 35 NA * SW 36-37.5: SAA 36 MW-4S-40 * SW 36-37.5: SAA 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 36 MW-4S-40 * SW 36-37.5': SAA 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 38 Wine Sand #2 Mine Sand #2 Mine Sand #2 | 20 | 2 | | * | SP | Reddish brown medium to coarse sand | l. wet | | | | |
| 25 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38 bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38 bgs 40 MW-4S-40 * SW Bio-5-37.51: SAA ST-5-381: Fine-medium grit brown sand 38-401: Medium to coarse sand SW Locking Cap Morie Sand #2 Morie Sand #2 Morie Sand #2 Morie Sand #2 | 20 | - | | | 51 | | ., | | | | |
| 25 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38 bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38 bgs 40 MW-4S-40 * SW Bio-5-37.51: SAA ST-5-381: Fine-medium grit brown sand 38-401: Medium to coarse sand SW Locking Cap Morie Sand #2 Morie Sand #2 Morie Sand #2 Morie Sand #2 | | | | | | | | | | | |
| 25 3.5 * SP SAA 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38 bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38 bgs 40 MW-4S-40 * SW Bio-5-37.51: SAA ST-5-381: Fine-medium grit brown sand 38-401: Medium to coarse sand SW Locking Cap Morie Sand #2 Morie Sand #2 Morie Sand #2 Morie Sand #2 | | | MW-48-23 5 | * | | Groundwater encountered at ~23 ft | | | | | |
| 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 40 N/A * SW 36.5-37.5': SAA 37.5-38': Fine-medium grit brown sand 37.5-38': Fine-medium grit brown sand 38-40': Medium to coarse sand Image: Subscript of the screened interval screened interval screened interval from 28-38' bgs Image: Subscript of the screened interval screened | | | 10100 10 20:0 | | | Groundwater encountered at 25 ft | | | | | |
| 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet Slotted screened interval from 28-38' bgs 40 N/A * SW 36.5-37.5': SAA 37.5-38': Fine-medium grit brown sand 37.5-38': Fine-medium grit brown sand 38-40': Medium to coarse sand Image: Subscript of the screened interval screened interval screened interval from 28-38' bgs Image: Subscript of the screened interval screened | 25 | 35 | | * | SP | SAA | | | | | |
| 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 40 N/A * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand Image: SP Image: SP S8-40': Medium to coarse sand Image: SP Image: SP <t< td=""><td>23</td><td>5.5</td><td></td><td></td><td>51</td><td>5. L L</td><td></td><td></td><td></td></t<> | 23 | 5.5 | | | 51 | 5. L L | | | | | |
| 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 40 N/A * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand Image: SP Image: SP S8-40': Medium to coarse sand Image: SP Image: SP <t< td=""><td></td><td></td><td></td><td> </td><td> </td><td> </td><td></td><td></td><td></td></t<> | | | | | | | | | | | |
| 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 40 N/A * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand Image: SP Image: SP SB-40': Medium to coarse sand Image: SP Image: SP <t< td=""><td></td><td></td><td></td><td></td><td></td><td> </td><td></td><td></td><td></td></t<> | | | | | | | | | | | |
| 30 MW-4S-30 * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 35 3.5 N/A * SW Brown medium-coarse sand, trace gravel, trace mica, wet from 28-38' bgs 40 N/A * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand Image: SP Image: SP SB-40': Medium to coarse sand Image: SP Image: SP <t< td=""><td></td><td></td><td></td><td></td><td></td><td> </td><td></td><td>Slotted screened interval</td><td></td></t<> | | | | | | | | Slotted screened interval | | | |
| 35 3.5 N/A * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38.40': Medium to coarse sand 40 MW-4S-40 * Morie Sand #2 | 30 | | MW-48-30 | * | SW | Brown medium-coarse sand trace gray | vel trace mica wet | | | | |
| 40 MW-4S-40 * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 40 MW-4S-40 * Morie Sand #2 Locking Cap Morie Sand #2 | 50 | | 00-01-11 | | 3 11 | biown meerani-coarse saile, trace gra | ver, trace inica, wet | 110111 20-00 Ugs | | | |
| 40 MW-4S-40 * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand 40 MW-4S-40 * SP 38-40': Medium to coarse sand Image: Second s | | | | | | | | | | | |
| 40 MW-4S-40 * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 40 MW-4S-40 * Morie Sand #2 Locking Cap Morie Sand #2 | | | | | | | | | | | |
| 40 MW-4S-40 * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit brown sand SP 38-40': Medium to coarse sand 40 MW-4S-40 * Morie Sand #2 Locking Cap Morie Sand #2 | | | | | | | | | | | |
| 40 MW-4S-40 * SW 36.5-37.5': SAA SM 37.5-38': Fine-medium grit browm sand SP 38-40': Medium to coarse sand 40 MW-4S-40 * Morie Sand #2 Locking Cap Morie Sand #2 | 25 | 25 | | * | | | | | | | |
| 40 MW-4S-40 * SM 37.5-38°: Fine-medium grit browm sand 40 MW-4S-40 * SP 38-40': Medium to coarse sand Locking Cap Morie Sand #2 | 35 | 3.3 | IN/A | Ľ | CW | 26 5 27 51, 8 4 4 | | | | | |
| 40 MW-4S-40 * SP 38-40': Medium to coarse sand Locking Cap Morie Sand #2 | | | | | | | . 1 | | | | |
| 40 MW-4S-40 * * | | | | | | | IU | | | | |
| Locking Cap Morie Sand #2 | | | | | SP | 56-40": Medium to coarse sand | | | | | |
| Locking Cap Morie Sand #2 | 40 | | MW AS 40 | * | | | | | | | |
| | 40 | | 1v1 vv -45-40 | ľ | | | | | | | |
| | | | I | I | I | I | | I | | | |
| Screened interval Bentonite seal | 1 | Locking | g Cap | | I | Morie Sand #2 | | | | | |
| | : | Screene | d interval | | 1 | Bentonite seal | | | | | |
| Grant | | | | | - | | | | | | |
| Grout Grout Outside temp. during drilling was too cold for the PID to work properly | : | * | Outside temp. duri | ng drillir | ng was to | | | | | | |

| | | | | | | Project: Cropsey | Boring ID MW-5 | | |
|-----------------|----------------|-------------------------------|-----------|----------------|--|--|---------------------------|----|--|
| | | | | | | Brooklyn, NY | Sheet 1 of 1 | | |
| | | tion: Grass nort | | ide of | Bay 25th | Project Manager: R. Costanzo | Project Number: 85265 | | |
| Ground | d Elev | ation: | 22.6 | | | Dated Started: 3/8/16 | Date completed: 3/9/16 | | |
| | | | | | | Drill Type: Direct Push for soil | | | |
| | | | | | | cores; Hollow stem auger for well | | | |
| Depth | to Firs | t Water: | 19 | | | installation | Borehole Dia: 4.25" | | |
| Depth | to Stat | tic Water: | 19.4 | | | Drill Rig Type: | Geoprobe | | |
| Top of | ` Casin | g Elevation: | 22.2 | | | Drilling Company: | Zebra | | |
| | | • | | Sampl | er | Driller's Name: | Luke | | |
| Size: 2 | " I.D. | PVC | Type: | | Encore | APEX Representative: | Ashlyn Norberg | | |
| Screen | : 0.010 |)-inch Slot PVC | | er: | NA | Owner/Client Rep.: | AMA Cropsey | | |
| Sand: # | #2 Mo | rie | Fall: | | NA | * | | | |
| | × | | (1 | | | L | | | |
| Depth (feet) | Recovery | Sample I.D. or Control No. | PID (ppm) | USCS Class. | Descri | ption of Soils | Well Construction | on | |
| 0 | N/A | N/A | 0 | SM | Topsoil/grass, grades to Silty sand, son | ne gravel, brown, moist | Locking Cap at grade | | |
| | | | | | Hand cleared | | | | |
| | | | | | | | | | |
| └ │ | | | | | | | | | |
| ا_ ا | NT/ 4 | NT/ A | NT/ A | | 1 | | | | |
| - 5 | N/A | N/A | N/A | | | | | | |
| - | | | | | | | | | |
| - | | | | | | | | | |
| - | | | | | | | | | |
| 10 | N/A | N/A | N/A | | | | | | |
| 10 | 1011 | 1011 | 1011 | | | | | | |
| - | | | | | | | | | |
| - | | | | | | | | | |
| - | | | | | | | | | |
| 15 | 3.5 | N/A | 0 | | 15-16': Brown medium grit sand, some | | | | |
| - | | | 0 | | 16-17': Medium to fine grit sand, light | | | | |
| - | | | 0 | SW | 17-18': Medium to fine grit sand, trace | | 1 | | |
| - | | | 0 | SW | 18-19': Brown medium grit sand, trace | gravel, dry; Groundwater encountered at appro- | ximately 19' bgs | | |
| | | | | ~ ~ | | | | | |
| 20 | 4 | | 0 | SP | 20-22.5': Medium to fine grit sand, red | aisn brown | Slattad compared internal | | |
| - | | MW-5-22.5 | 100 | CM/ | 22.5.25' Madium cuit and tung | al wat | Slotted screened interval | | |
| - | | IVI VV -3-22.3 | 100 | SW | 22.5-25': Medium grit sand, trace grave | cı, wei | from 17' - 27' bgs | | |
| - | | | | | | | | | |
| 25 | 4 | N/A | | SW | 25-29': SAA | | | | |
| | | | | | | | | | |
| - | | | | | | | | | |
| | | | | | | | | | |
| _ | | | | | | | | | |
| 30 | 4.8 | MW-5-30 | 131 | SW | SAA | | | | |
| - | | | | | 1 | | | | |
| - | | | | | | | | | |
| - | | | | | | | | | |
| 35 | | MW-5-35 | 101 | SW | SAA | | | | |
| 55 | | | 101 | 5.1 | | | 1I | | |
| | Locking | g Cap | | | Morie Sand #2 | | | | |
| | Screene | ed interval | | | Bentonite seal | | | | |
| | | | | | | | | | |
| | | | | | Grout | | | | |
| | | | | | | | | | |

| | | | | | | Project: Cropsey | Boring ID MV | W-5I |
|-----------------|----------|------------------|-----------|----------------|---|--|---------------------------|---------|
| | | | | | | Brooklyn, NY | Sheet 1 of 1 | |
| Bori | ngIc | cation: Sidewalk | on the | corner | of Cropsey Ave and 20th Ave | Project Manager: R. Costanzo | Project Number: 85265 | 5 |
| Grou | ind F | levation: | on the | conner | of cropsey rive and 20th rive | Dated Started: | Date completed: 6/23/ | |
| GIU | ind L | evation. | | | | Drill Type: Direct Push for soil cores; Hollow | Date completed. 0/25/ | 10 |
| Dent | 1. 4 . T | | | | | | Borehole Dia: 4.25" | |
| | | First Water: | | | | stem auger for well installation | | |
| Dept | h to S | Static Water: | | | | Drill Rig Type: | Geoprobe | |
| Тор | of Ca | sing Elevation: | - | | | Drilling Company: | Zebra | |
| | | | | Samp | | Driller's Name: | Luke | |
| Size: | 2" I. | D. PVC | Type: | | Encore | APEX Representative: | Dan Haug | |
| Scree | en: 0. | 010-inch Slot PV | Hamn | ner: | NA | Owner/Client Rep.: | AMA Cropsey | |
| Sand | l: #2 I | Morie | Fall: | | NA | | | |
| | | ۲ | (u | | | | | |
| Depth (feet) | Recovery | Sample I.D. of | bu | USCS Class. | _ | | | |
|)ep | | Control No. | D (F | JSC Class | D | Description of Soils | Well Const | ruction |
| U C | - Be | | PID (ppm) | 50 | | | | |
| | 0 N/ | A N/A | | | Topsoil/grass, grades to Silty sand, so | me gravel brown moist | Locking Cap at grade | |
| - | 0 10/ | AIWA | 0 | Í l | Hand cleared | ne gravel, brown, moist | Locking Cap at grade | |
| - | | | | | Tiand cleared | | | |
| - | | | | | | | | |
| F | | | | | | | | |
| F | 5 N/ | A N/A | N/A | | | | | |
| - | - | | | | | | | |
| F | | | | | | | | |
| Γ | | | 1 | 1 | | | | |
| - | | | | | | | | |
| 1 | 0 N/ | A N/A | N/A | | | | | |
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| _ | | | | | | | | |
| 1 | 5 N/A | N/A | N/A | | | | | |
| - | | | | | | | | |
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| 2 | 20 | | | | | | | |
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| _ 2 | 25 | | | | | | | |
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| L | | | 1 | 1 | | | Slotted screened interval | |
| 3 | 30 | | | | | | from 40-50' bgs | |
| L | | | 1 | 1 | | | | |
| L | | | | | | | | |
| L | | | | 1 | | | | |
| L | | | | 1 | | | | |
| 3 | 5 | | | | | | | |
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| L | | | | | | | | |
| 4 | 10 | | | | | | | |
| L | | | | | | | | |
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| | 15 | | | | | | | |
| 4 | 15 | | | | | | | |
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| | | | | | | | Project: Cropsey | Boring ID MW | 7-6 | |
|-----------------|----------|-------------------------------|-----------|-------|----------------------------------|-------------|--------------------------------------|---------------------------------|---------|--|
| | | | | | | | Brooklyn, NY | Sheet 1 of 1 | | |
| | | tion: Grass sout | | | f Bay 25th | | Project Manager: R. Cosntanzo | Project Number: 85265 | | |
| Groun | d Eleva | ation: | 23.2 | 2 | | | Dated Started: 3/8/2016 | Date completed: 3/9/16 | | |
| | | | | | | | Drill Type: Direct Push for soil | | | |
| | · | | 27 | - | | | cores; Hollow stem auger for well | D 1 1 D' 4 25" | | |
| | | t Water: | 27 | | | | installation | Borehole Dia: 4.25" Geoprobe | | |
| | | ic Water: g Elevation: | 20. | | | | Drill Rig Type: | Zebra | | |
| 100 01 | Casin | g Elevation: | 22.3 | | pler | | Drilling Company: Driller's Name: | Zebra Luke | | |
| Size: 2 | "חו | PVC | Туре | | Encore | | APEX Representative: | Ashlyn Norberg | | |
| | |)-inch Slot PVC | | | NA | | Owner/Client Rep.: | AMA Cropsey | | |
| Sand: 7 | | | Fall: | iner. | NA | | owner/enent rep. | Thin Cropsey | | |
| | | | | T | | | | | | |
| Depth (feet) | Recovery | Sample I.D. or Control No. | PID (ppm) | USCS | Class. | Descri | ption of Soils | Well Constr | ruction | |
| 0 | N/A | N/A | | 0 SN | 1 Topsoil/grass, grades to Silty | sand, son | ne gravel, brown, moist | Locking Cap at grade | | |
| [| | | | | Hand cleared | | | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| - 5 | N/A | N/A | N/A | | | | | | | |
| - | | | | | | | | | | |
| _ | | | | | | | | | | |
| - | | | | | | | | | | |
| 10 | N/A | N/A | N/A | | | | | | | |
| 10 | IN/A | 11/74 | IN/A | | | | | | | |
| - | | | | | | | | | | |
| _ | | | | | | | | | | |
| | | | | | | | | | | |
| _ 15 | 4 | N/A | | 0 G | 15-19': Light brown medium | grit sand, | course gravel, trace mica, dry | | | |
| - | | | | | | | | | | |
| - | | | | | | | | | | |
| | | | | | | | | | | |
| 20 | 5 | MW-6-23 | | 0 G | | | | | | |
| - | | | | 0 SW | 21.5-25': Reddish brown med | dium grit s | and, trace gravel | Slotted screened interval | | |
| - | | | | | | | | from 17' - 27' bgs | | |
| - | | | | | | | | | | |
| - 25 | 5 | | | 0 SV | 25-27 Dark brown mediun gr | | | | | |
| | | | | SI | 27-29': Reddish brown mediu | um grit sai | nd, wet | | | |
| - | | | | | Groundwater encountered at | approxim | ately 27' bgs | | | |
| - | | | | | | | | | | |
| - 30 | | MW-6-30 | , | 0 51 | Reddish brown medium grit s | sand, wet | | | | |
| | | | | | ere with moutain gift a | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| 35 | | 1 | | | <u> </u> | | | <u>I</u> | | |
| | Locking | g Cap | | | Morie Sand #2 | | | | | |
| | Screene | d interval | | | Bentonite seal | | | | | |
| | Sercent | a morvar | | | Demonite Scal | | | | | |
| | | | | | Grout | | | | | |
| | | | | | | | | | | |

| | | | | | | Project: Cropsey | Boring ID MW | 7 | | |
|-----------------|----------|-------------------------------|-----------|----------------------|--|---|--|--------|--|--|
| | | | | | | Brooklyn, NY | Sheet 1 of 1 | -/ | | |
| Boring | T Locat | tion: Grass sout | heast c | orner (| of Bay 25th and Cropsey Ave | Project Manager: R. Costanzo | Project Number: 85265 | | | |
| Groun | | | 23.5 | | or Day 25th and Cropsey Ave | Dated Started: 3/8/2016 | Date completed: 3/8/16 | | | |
| Oloui | u Elev | ation. | 23.3 | | | Drill Type: Direct Push for soil cores; Hollow | | | | |
| Donth | to Fire | t Water: | 23 | | | stem auger for well installation Borehole Dia: 4.25" | | | | |
| | | ic Water: | 20.1 | | | | rill Rig Type: Geoprobe | | | |
| | | g Elevation: | 20.1 | | | Drilling Company: | Zebra | | | |
| 1000 | Casili | g Elevation. | 23.1 | Samp | lor | Driller's Name: | Luke | | | |
| Sizo. | 2" I.D. | DVC | Type: | - | Encore | APEX Representative: | Ashlyn Norberg | | | |
| | |)-inch Slot PVC | Hamn | oer: | NA | Owner/Client Rep.: | AMA Cropsey | | | |
| | #2 Mo | | Fall: | ICI. | NA | Owner/Chent Rep | AWA Cropsey | | | |
| Sanu. | r. | | | | NA | | 1 | | | |
| Depth (feet) | Recovery | Sample I.D. or Control No. | PID (ppm) | USCS Class. | Γ | Description of Soils | Well Constru | uction | | |
| 0 | N/A | N/A | 0 | SM | Topsoil/grass, grades to Silty sand, so Hand cleared | me gravel, brown, moist | Locking Cap at grade | | | |
| 5 | N/A | N/A | N/A | | | | | | | |
| 10 | N/A | N/A | N/A | | | | | | | |
| 15 | 3 | N/A | 0 | GW SW GP SW | 15-15.5': Coarse gravel 15.5-16': Medium grit brown sand 16-17': Coarse gravel, medium grit sa 17-18':Medium grit grey/brown sand, | | | | | |
| 20 | 4 | MW-7-23 | 0 | SW SW SM SW | 20-21.5': Dark brown medium grit sar 21.5-23': Grey medium grit sand 23-23.5': Brown fine sand, wet; Groun 23.5-24': Brown medium grit sand | nd, dry ndwater encountered at approximately 23' bgs | Slotted screened interval from 20-30' bgs | | | |
| 25 | 4 | N/A | 0 | sw | 25-29': Reddish brown medium grit sa | and, trace gravel | | | | |
| 30 | 4.8 | 5 | 0 | sw | SAA | | | | | |
| 35 | | MW-7-34 | | | | | | | | |
| | Locking | g Cap | | I | Morie Sand #2 | | | | | |
| | Screene | d interval | | | Bentonite seal | | | | | |
| | | | | | Grout | | | | | |

| Biooklyn, NY Shee 1 of 1 Brondy Learning Shee 1 of 1 Ground Elevation: Tops Charger R. Costamov Project Number: 8526 Depth b First Water: Data Startici: 6242016 Depth b First Water: Booklyn, NY Depth b First Water: Booklyn, NY Depth b First Water: Booklyn, NY Top of Casing Elevation: Booklyn, NY Size: 2° LD, PVC Type: Size: 2° LD, PVC Type: Booklyn, Starting Company: Size: 2° LD, PVC Type: Booklyn, NY Starting Size: 2° LD, PVC Type: Booklyn, NY Starting Size: 2° LD, PVC Type: Booklyn, NY Company: Size: 2° LD, PVC Type: Booklyn, NY Starting Size: 2° LD, PVC Starting <th></th> <th></th> <th></th> <th></th> <th></th> <th></th> <th>Project: Cropsey</th> <th>Boring ID MV</th> <th>V-8</th> | | | | | | | Project: Cropsey | Boring ID MV | V-8 | |
|--|----------------------------------|--------------|-----------------|-----------|----------------|---------------------------------------|--|-------------------------|-----|--|
| Boring Location: Grass, Southwest Courtyard Project Manager, R. Costanzo Project Manager, R. Costanzo Ground Elevation: Dated Statel: 624/016 Date complete: 624/16 Depth to Static Water: 18.3 Depth to Static Water: 18.3 Dig to Grassing Elevation: Dill Type: Direct Push for soil cores; Hollow Borehold: Dia: 4.25" Depth to Static Water: 12.2 Dill Type: Direct Manager, R. Costanzo State: 2" LD. PVC Type: Encore State: 2" Advise Date State: Date State: State: 2" Advise: Date State: Date State: State: 2" Advise: Fall: NA O NA NA 0 State: 2" State: State: State: State: 2" State: State: State: State: State: State: State: State: State: </td <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>-</td> <td></td> | | | | | | | | - | | |
| | Boring | Locat | ion: Grass, Sou | thwest | Court | vard | | | 5 | |
| $ \begin{array}{c c c c c c c c c c c c c c c c c c c $ | | | | | | | | | | |
| Depth to Static Water: 18.3 Delil Rig Type: Geoprobe Top of Casing Elevation: Samplet Driller's Name: Luke Size: 2" LD. PVC Type: Encore APEX Representative: Dan Kopce Size: 2" LD. PVC Type: Encore APEX Representative: Dan Kopce Sand: #2 Morie Fall: NA Owner/Client Rep: AMA Cropsey Sand: #2 Morie Fall: NA Owner/Client Rep: AMA Cropsey Sand: #2 Morie Fall: NA Owner/Client Rep: Owner/Client Rep: Sand: #2 Morie Fall: NA Owner/Client Rep: Owner/Client Rep: Sand: #2 Morie Fall: NA NA Owner/Client Rep: Owner/Client Rep: Sand: #2 Morie Sample LD. or Sample Sample LD. or Sample S | | | | | | | | w | | |
| Depth to Static Wate: 18.3 Drill Rig Type: Geoprobe Top of Casing Heration: Sing 2: 10 PVC Type: Encore APEX Representative: Dan Kopee Size: 2" LO PVC Type: Encore APEX Representative: Dan Kopee Size: 2" Morie Fall: NA Owner/Client Rep: AMA Cropsy Sand: 72 Morie Fall: NA Owner/Client Rep: Mark Copsy Sand: 72 Morie Fall: NA Owner/Client Rep: Mark Copsy Sand: 72 Morie Fall: NA Owner/Client Rep: Mark Copsy Sand: 72 Morie Fall: NA Owner/Client Rep: Description of Soils Well Construction 10 NA NA NA NA SW 9 - 16" Tan/Dark Brown Fine Sand, Moist Lacking Cap at grade 11 NA NA SW 9 - 16" Tan/Dark Brown Fine Sand, Moist GW encountered at 19.5" 12 NA NA SW 9 - 16" Tan/Dark Brown Fine Sand, Moist GW encountered at 19.5" 13 NA NA SW 9 - 16" Tan/Dark Brown Fine Sand, Moist GW encountered at 19.5" < | Depth | to Firs | t Water: | | | | stem auger for well installation | Borehole Dia: 4.25" | | |
| Junc Sampler Driller's Marrie' Luke Strict 2" LD PVC Type: Encore $APEX$ Representative: Dan Kopee Serven: 0.010-inch Slot PVC Harmor: NA Owner/Clean Rep.: AMA Cropsey Sand: #2 Morie Fall: NA Owner/Clean Rep.: AMA Cropsey $g \in g : g : g : g : g : g : g : g : g : $ | | | | 18.3 | | | | | | |
| Size: 2*1 LD PVC Type: Encore APEX Representative: Dan Kopee Server: 0:010-inch Slot PVC Hammer. NA Owner/Client Rep.: AMA Cropsey Sand: #2 Morie Fall: NA $\frac{g}{g}$ < | Top of | Casin | g Elevation: | | | | Drilling Company: | | | |
| Server: 0.01-inch Slov PVC Harmer: NA Owner/Client Rep.: AMA Cropsey Sand: #2 Noric Fall: NA Vector | | | | | Sampl | er | Driller's Name: | | | |
| Sand: $\#2$ Moric Fall: NA $\frac{6}{20}$ $\frac{6}{20}$ $\frac{6}{20}$ $\frac{7}{20}$ $\frac{7}{20}$ Description of Soils Well Construction $\frac{6}{20}$ $\frac{6}{20}$ $\frac{7}{20}$ $\frac{7}{20}$ $\frac{7}{20}$ $\frac{7}{20}$ Description of Soils Well Construction 0 NA NA 0 SM Topoligness grades to Sily sand, some gravel, brown, moist Locking Cap at grade 10 0.5 NA NA 0 SW $9.5 \cdot 10^\circ$: Tan/Dark Brown Fine Sand, Moist Locking Cap at grade 10 0.5 NA 0 SW $9.5 \cdot 10^\circ$: Tan/Dark Brown Fine Sand, Moist Locking Cap at grade 10 0.5 NA 0 SW $9.5 \cdot 10^\circ$: Tan/Dark Brown Fine Sand, Moist Control Tan/Dark Brown Mediam-Coarse Sand, Little - no fines Common Sand, Moist Control Tan/Dark Brown Fine Sand, Moi | | | | Type: | | | | | | |
| End Description of Soils Well Construction 0 N/A N/A 0 SM Topsol/grass, grades to Sily sand, some gravel, brown, moist Locking Cap at grade 10 0.5 N/A N/A 0 SW 9.5 - 10° : Tan/Dark Brown Fine Sand, Moist 10 0.5 N/A 0 SW 9.5 - 10° : Tan/Dark Brown Fine Sand, Moist 11 0.5 N/A 0 SW 12 - 12.5° : Tan Fine Sand, Moist 115 3 N/A 0 SW 12 - 12.5° : Tan Fine Sand, Moist 12 1.2 1.2.8° : Brown Orange Clay, Moist 0 SW 13 3 N/A 0 SW 12 - 12.5° : Tan Fine Sand, Moist 14 0.5 N/A 0 SW 12 - 12.8° : Brown Orange Clay, Moist 15 3 N/A 0 SW 12 - 12.8° : Brown Modium-Coarse Sand, Linte- no fines (0.5%) 17 1.7 1.7 1.7° : Gray Clay, Wet 20 20 3 0 SW 18 - 20: Brown Medium-Coarse Sand, Linte- no fines 20.5° : 23.3° : Brown Fine Sand, Wet 21 4.8 MW-8 (307) 0 SW 25 - 27.4°: Dark Brown Fine Sand, Wet 30 4.8 MW-8 (307) 0 SW 25 | | | | | er: | | Owner/Client Rep.: | AMA Cropsey | | |
| 0 N/A N/A 0 SM Topsoil/grass, grades to Silty sand, some gravel, brown, moist Looking Cap at grade 5 N/A N/A 0 SW 9.5 - 10': Tan/Dark Brown Fine Sand, Moist 10 0.5 N/A 0 SW 9.5 - 10': Tan/Dark Brown Fine Sand, Moist 11 0.5 N/A 0 SW 12.12.5: Tan Fine Sand, Moist 0 SW 12.12.5: Tan Fine Sand, Moist 0 SW 15 3 N/A 0 GH 13.7 - 15': Brown Cores Sand, Little - no fines (0.5%) 17 17.4 17.7: roary Clay, Wet 13.7 - 15': Brown Medium-Coarse Sand, Moist GW encountered at 19.5' 20 3 0 GH 23.3' - 25': Gmy Clay, Wet Slotted screened interval from 17.27 bgs 21 4.8 MW-8 (30') GW 25' - 27.4': Dark Brown Fine-Medium-Coarse Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand, Wet Slotted screened interval from 17.27 bgs 23 4.8 MW-8 (30') GW 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand, Wet Slotted screened interval from 17.27 bgs 35 < | Sand: # | #2 Mo | rie | Fall: | | NA | | | | |
| 5N/AN/A0100.5N/A0110.5N/A0120SW9.5 - 10 : Tan/Dark Brown Fine Sand, Moist130SW12 - 12.5 : Tan Fine Sand, Moist140SW12 - 12.5 : Tan Fine Sand, Moist153N/A0SW153N/A0160CH13.2 - 13.7 : Gray Clay, Wet1718717.1 7.7 : Brown Caraes Sand, Moist180CH17.4 - 17.7 : Gray Clay, Wet193W18' - 20': Brown Medium-Coarse Sand, Little - no fines2030CH17.7 - 187: Tan Fine Sand, MoistGW encountered at 19.5'2030CH212.5 - 23.3: Brown Fine-Medium Sand, WetSlotted screened interval from 17-27 bgs224.5N/A0SW2525.5 - 27.4: Dark Brown Fine Sand, Wet0264.8MW-8 (30)W274.8MW-8 (30)GW284.8MW-8 (30)0294.8MW-8 (30)0205.814.42115.815.8224.5NA231.81.8241.81.1251.81.8261.81.1271.81.1281.81.8291.81.8201.81.82 | Depth (feet) | Recovery | - | PID (ppm) | USCS Class. | D | Description of Soils Well Construction | | | |
| 10 0.5 N/A 0 SW 12 · 12.5 : Tan Fine Sand, Moist 15 3 N/A 0 CL 12.5 · 12.8 : Brown/Orange Clay, Moist 15 3 N/A 0 CH 13.7 · 132 : Tan Fine Sand, Moist 16 0 SW 12.7 · 132 : Tan Fine Sand, Moist 0 17 17 17 · 17 · 139 : Drown Medium-Coarse Sand. Little - no fines (0-5%) 0 17 · 17 · 179 : Drown Medium-Coarse Sand. Little - no fines (0-5%) 20 3 MW-8 (19.5) 0 CH 17.7 · 189 : Tan Fine Sand, Moist GW encountered at 19.5' 20 3 W/A 0 CH 17.7 · 179 : Grow Charse Sand. Little - no fines 20.5 · 23.3': Brown Medium-Coarse Sand. Little - no fines 20.5 · 23.3': Brown Fine-Medium Sand, Wet Slotted screened interval from 17-27' bgs 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 0 30 4.8 MW-8 (30') 0 SW 25' - 30': Black Gravelly Sand. Little-no fines. Wet 0 35 0 0 0 0 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet 0 | - - - | Hand cleared | | | | | ne gravel, brown, moist | Locking Cap at grade | | |
| 15 3 N/A 0 SW 12.75' - 13.2': Tan Fine Sand, Moist 15 3 N/A 0 CH 13.2 - 13.7': Gray Clay, Wet 0 SW 13.7 - 15': Brown Medium-Coarse Sand, Moist 6W encountered at 19.5' 0 HW-8 (19.5') 0 CH 17.4': Brown Coarse Sand, Moist 20 3 0 SW 12.75' - 13.2': Tan Fine Sand, Moist GW encountered at 19.5' 20 3 0 CH 17.4' - 17.7': Gray Clay, Wet GW encountered at 19.5' 20 3 0 SW 18' - 20': Brown Medium-Coarse Sand, Little - no fines 20.5' - 23.3': Brown Fine-Medium Sand, Wet Slotted screened interval from 17-27' bgs 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 0 30 4.8 MW-8 (30') 0 Slotted screened interval from 17-27' bgs 16W encountered at 19.5' 35 0 4.8 MW-8 (30') 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 0 36 0 0 0 0 0 0 0 0 0 0 | 10 | 0.5 | N/A | 0 | SW | 12 - 12.5' : Tan Fine Sand, Moist | | | | |
| 15 3 N/A 0 CH 13.2 - 13.7 : Gray Clay, Wet 0 5W 13.7 : 15 : Brown Medium-Coarse Sand. Little- no fines (0-5%) 17 - 17.4: Brown Coarse Sand, Moist GW encountered at 19.5' 20 3 0 CH 17.4 : 17.7: Gray Clay, Wet GW encountered at 19.5' 20 3 0 CH 17.4 : 17.7: Gray Clay, Wet GW encountered at 19.5' 20 3 0 CH 17.7 : 18: Tan Fine Sand, Moist GW encountered at 19.5' 20 3 0 CH 17.7 : 18: Tan Fine Sand, Moist GW encountered at 19.5' 20 3 0 CH 12.3 : - 25: Gray Clay, Wet Slotted screened interval from 17-27' bgs 21 4.5 N/A 0 SW 25' - 27.4: Dark Brown Fine Sand, Wet 27.4 - 28.5: Brown Medium-Coarse Sand, Wet 30 4.8 MW-8 (30') 0 GW 28.5' - 30: Black Gravelly Sand. Little- no fines. Wet 35 0 4.8 MW-8 (30') 0 0 0 35 0 0 0 0 0 0 0 | - | | | | | | pist | | | |
| 20 3 0 SW 13.7 · 15' : Brown Medium-Coarse Sand. Little- no fines (0-5%) 20 3 0 CH 17.4 · 17.7 · Gray Clay, Wet 20 3 0 SW 18' · 20': Brown Medium-Coarse Sand. Little - no fines 20 3 0 CH 17.4 · 17.7 · 17.8 : Tan Fine Sand, Moist GW encountered at 19.5' 20 3 0 SW 18' · 20': Brown Medium-Coarse Sand. Little - no fines Slotted screened interval from 17 · 27' bgs 20 3 0 CH 23.3' · 25': Gray Clay, Wet Slotted screened interval from 17 · 27' bgs 25 4.5 N/A 0 SW 25' · 27.4': Dark Brown Fine Sand, Wet 27.4' · 28.5': Brown Medium-Coarse Sand, Wet 25. · 30': Black Gravelly Sand. Little-no fines. Wet 10 30 4.8 MW-8 (30') 0 GW 28.5' · 30': Black Gravelly Sand. Little-no fines. Wet 10 35 10 10 10 10 10 10 | 1.5 | 2 | NT/ A | | | - | | | | |
| 20 3 0 CH 17 - 17.4': Brown Coarse Sand, Moist 20 3 0 CH 17.4': -17.7': Gray Clay, Wet GW encountered at 19.5' 20 3 0 SW 18': 20': Brown Medium-Coarse Sand. Little - no fines Slotted screened interval 20 3 0 CH 23.3': 25': Gray Clay, Wet Slotted screened interval 20 - - - - - - 21 - - CH 23.3': 25': Gray Clay, Wet Slotted screened interval 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet Slotted screened interval 30 4.8 MW-8 (30') GW 25' - 27.4': Dark Brown Fine Sand, Wet 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet Interval 35 Interval Interval Interval Interval Interval 36 Interval Interval Interval Interval Interval 37 Interval Interval Interval Interval Interval 38 Interval Interval Interval Inter | - 15 | 3 | 1N/A | | | | nd Little- no fines (0-5%) | | | |
| 20 3 MW-8 (19.5') 0 CH 17.4' - 17.7': Gray Clay, Wet GW encountered at 19.5' 20 3 0 SW 18' - 20': Brown Medium-Coarse Sand. Little - no fines Slotted screened interval 20 3 0 CH 2.3' - 25': Gray Clay, Wet Slotted screened interval 21 0 CH 2.3.3' - 25': Gray Clay, Wet Slotted screened interval 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand. Wet 30 4.8 MW-8 (30') GW 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet Image: Comparison of the stand of the s | - | | | | 5.0 | | | | | |
| 20 3 0 SW 18' - 20': Brown Medium-Coarse Sand. Little - no fines 20.5' - 23.3': Brown Fine-Medium Sand, Wet Slotted screened interval from 17-27' bgs 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet Slotted screened interval from 17-27' bgs 26 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet 30 4.8 MW-8 (30') GW 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet Image: Coarse Sand, Wet Image: Coarse Sa | - | | MW-8 (19.5') | 0 | CH | | | | | |
| 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet Slotted screened interval from 17-27' bgs 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 10 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 10 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet 10 1 | _ | | | 0 | | | | GW encountered at 19.5' | | |
| 25 4.5 N/A 0 CH 23.3' - 25': Gray Clay, Wet 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 0 GW 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 30 4.8 MW-8 (30') GW 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet 35 | 20 | 3 | | 0 | SW | | | | | |
| 25 4.5 N/A 0 SW 25' - 27.4': Dark Brown Fine Sand, Wet 0 0 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 30 4.8 MW-8 (30') 0 GW 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet 35 | - | | | 0 | CH | 23.3' - 25': Gray Clay, Wet | | 10111727 053 | | |
| 30 4.8 MW-8 (30') 0 27.4' - 28.5': Brown Medium-Coarse Sand, Wet 30 4.8 MW-8 (30') 0 GW 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet 35 | - | | | | | | | | | |
| 30 4.8 MW-8 (30') 0 GW 28.5' - 30': Black Gravelly Sand. Little-no fines. Wet 35 | 25 | 4.5 | N/A | 0 | SW | 25' - 27.4': Dark Brown Fine Sand, We | et | | | |
| 30 4.8 MW-8 (30') 35 | | | | | GW | | | | | |
| 35 | - | | | 0 | 0.11 | 20.0 50. Black Graveny Sand. Effic | | | | |
| | 30 | 4.8 | MW-8 (30') | | | | | | | |
| | _ | | | | | | | | | |
| | - | | | | | | | | | |
| | - | | | | | | | | | |
| | 35 | | | | | | | | | |
| | | Locking | g Cap | | | Morie Sand #2 | | l | | |
| Screened interval Bentonite seal | Screened interval Bentonite seal | | | | | Bentonite seal | | | | |
| | | | | | | | | | | |
| Soil Cuttings Fill Grout | | Soil Cu | tings Fill | | | Grout | | | | |

| | | | | | | Project Company | Boring ID MV | 7.0 |
|---|----------|-------------------------------|--------------------------------------|----------------|--|---|---|-----|
| | | | | | | Project: Cropsey | 0 | V-9 |
| Devine Leasting Cross Southwart C. (1 | | | | | | Brooklyn, NY | Sheet 1 of 1 Project Number: 85265 | |
| Boring Location: Grass, Southwest Courtyard Ground Elevation: | | | | | | Project Manager: R. Costanzo | | |
| | | | | | | Dated Started: 6/24/2016 | Date completed: 6/23/1 | 0 |
| | | | | | | Drill Type: Direct Push for soil cores; Hollow stem auger for well installation | Parahala Dia: 4.25" | |
| | | | | | | | Borehole Dia: 4.25" | |
| | | | | | | Drill Rig Type: | Geoprobe Zebra | |
| | | | | | | Drilling Company: | Luke | |
| | | | | | | Driller's Name: | | |
| Size: 2" I.D. PVC Type: Encore Screen: 0.010-inch Slot PVC Hammer: NA | | | | | Encore | APEX Representative: | Dan Haug | |
| Screen: 0.010-inch Slot PVC Sand: #2 Morie | | | Fall: NA | | | Owner/Client Rep.: | AMA Cropsey | |
| Sand: # | he | | | | | | | |
| Depth (feet) | Recovery | Sample I.D. or Control No. | PID (ppm) | USCS Class. | Description of Soils Well Construction | | ruction | |
| - 0 | N/A | N/A | 0 | SM | Topsoil/grass, grades to Silty sand, sor Hand cleared | ne gravel, brown, moist | Locking Cap at grade | |
| 5 10 | N/A 2 | N/A N/A | N/A 0 10.7 23.4 0 0.4 | CL SW | 8' - 9.7': Brown Medium Sand. Trace g 9.7' - 10': Light Tan Fine Sand. Trace 1 11.3' - 11.7': Light Tan Fine Sand. Slig 11.7' - 13': Dark Gray Clay. Wet 13' - 22': Brown Medium-Coarse Sand | | | |
| 15 | 3.7 | N/A MW-9 (19.4') | 0 | | | | | |
| 20 | 3.7 | | 0 0 | | 22' - 24': Dark Brown/Gray Fine-Medi 24' - 25': Dark Gray Clay. Wet | um Sand. Wet | GW encountered at 19.4' Slotted screened interval from 17-27' bgs | |
| 25 | 4 | | | | 26.4' - 27.4': Dark Brown Medium-Co 27.4' - 28': Dark Brown Gravel. Little- 28' - 28.7': Brown Gravel. Some sand (28.7' - 30': Coarse Sand. Wet | no fines. | | |
| 30 | 3.7 | MW-9 (30') | | | 20.7 - 50 . Coaise Sailti. Wet | | | |
| 35 | | | | | | | | |
| Locking Cap Morie Sand #2 | | | | | Morie Sand #2 | | 1 | |
| Screened interval | | | | | Bentonite seal | | | |
| Soil Cuttings Fill | | | | | Grout | | | |

| | | | | | | Durainati Changay | Doming ID MI | W-10 | |
|-----------------|------------|-------------------------------|---------------|----------------|---|---|---|----------|--|
| | | | | | | Project: Cropsey | 0 | w-10 | |
| р [.] | т., | · | .1 | <u> </u> | 1 | Brooklyn, NY | Sheet 1 of 1 Project Number: 8526 | 5 | |
| Boring | Locat | ion: Grass, Sou | itnwest | Court | yard | Project Manager: R. Costanzo | | | |
| Ground | a Eleva | ation: | | | | Dated Started: 6/24/2016 | Date completed: 6/24/ | 16 | |
| | | | | | | Drill Type: Direct Push for soil cores; Hollow | D 1 1 D' 405" | | |
| | | t Water: | 10.0 | | | stem auger for well installation | Borehole Dia: 4.25" | | |
| | | ic Water: | 19.3 | | | Drill Rig Type: | Geoprobe | | |
| Top of | Casin | g Elevation: | | ~ 1 | | Drilling Company: | Zebra | | |
| ~ • | | | - | <u>Sampl</u> | | Driller's Name: | Luke | | |
| Size: 2 | | | Type: | | Encore | APEX Representative: | Dan Kopec | | |
| | |)-inch Slot PVC | | ner: | NA | Owner/Client Rep.: | AMA Cropsey | | |
| Sand: # | | rie | Fall: | | NA | | | | |
| Depth (feet) | Recovery | Sample I.D. or Control No. | PID (ppm) | USCS Class. | D | escription of Soils | Well Cons | truction | |
| - 5 | N/A N/A | N/A N/A | 0 N/A 0 | SM SW | Topsoil/grass, grades to Silty sand, son Hand cleared 6' - 8.4': Dark Brown Fine-Medium Sa 8.4' - 10': White, Dry, Concrete | | Locking Cap at grade | | |
| 10 | 4 | N/A | N/A 0 0 | SW | 12.5' - 13.1': White, Dry, Concrete 13.1' - 15': Light Brown Medium Sand | . Little-no fines (0-5%) | | | |
| 15 | 2.5 | N/A | 0 0 0 | SM SP SW | 16' - 17': Brown Fine Sand. Moist 17' - 17.5': Light Gray Fine Sand w/ Gi 17.5' - 20': Brown Medium-Coarse Sar | | | | |
| 20 | 4 | MW-10 (19.5') | 0 | SW | 21' - 21.8': Brown Medium-Coarse Sar 21.8' - 23.5': Light Brown Fine-Mediu 23.5' - 23.7': Dark Brown Fine Sand. S | m Sand. Wet | GW encountered at 19.5' Slotted screened interval from 17-27' bgs | | |
| 25 | 4 | N/A | 0 | | 23.7' - 25': Dark Brown Medium-Coars 25' - 26': - Light Brown Medium-Coars 26' - 29.1': Light Brown Fine-Medium 29.1' - 30': Dark Brown Medium-Coars | se Sand. Trace Gravels se Sand. Wet Sand. Wet | | | |
| 30 | 5 | MW-10 (30') | 0 | | | | | | |
| | | | | | | | | | |
| 35 | | | | | | | | | |
| 1 | Locking | , Cap | | l | Morie Sand #2 | | | | |
| : | Screene | d interval | |] | Bentonite seal | | | | |
| : | Soil Cut | tings Fill | |] | Grout | | | | |

| CLIENT: | WELL ID: MW - 15 |
|---------------------------|---|
| PROJECT #: (+0psly | ADDRESS: 2002-2004 Cropay Are |
| DATE: 2/10/16 | SAMPLING CREW: Dan + Ashlyn |
| DEPTH TO WATER (DTW) | 19.63 St DTB=22.0 ft |
| CASING DIAMETER (inches): | $\frac{\sqrt{2}}{3}$ <u>4</u> <u>4.5</u> <u>5</u> <u>6</u> <u>8</u> <u>12</u> Other |

PRODUCT PRESENT: _____ PRODUCT THICKNESS: _____

| TIME (2400 Hr) | VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE °F/°C | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
|-------------------|------------------|---------------|-------------------------|----------------------|---------------|-----|--------------------|-------------------|
| 1350 | 16 | 6.59 | 0.703 | 19.76 | 5,84 | 158 | 178 | Murky/Gray |
| 1355 | 18 | 6.56 | 0.729 | 21.56 | 4.88 | 146 | 87.3 | Chear |
| 1400 | 19 | 6.52 | 0.746 | 20.95 | 3.24 | 130 | 114 | Murky |
| 1405 | 20 | 6.50 | 0.743 | 21.53 | 3.27 | 121 | 82.7 | Clear |
| 1410 | 20.8 | 6.47 | 0,743 | 21.60 | 3.41 | 112 | 49.1 | Chear |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):____

| PURGING E | QUIPMENT | SAMPLING I | EQUIPMENT |
|---------------------|--------------------------|---------------------|--------------------------|
| 2" Submersible Pump | Bailer (Teflon®) | 2" Submersible Pump | Bailer (Teflon®) |
| Centrifugal Pump | Bailer (PVC) | DDL Sampler | Bailer (Stainless Steel) |
| 4" Submersible Pump | Bailer (Stainless Steel) | | Peristaltic Pump |
| Dedicated | Peristaltic Pump | Disposable Bailer | Dedicated Bailer |
| er: | 52 | Other: | |

WALL CONDITION: _____ GROUT: _____ MANWAY: _____ LOCKING CAP: _____ LOCK: _____ REMARKS: Mo-iba -21780 Monsoon 6W Pump - 16773

COMPLETED BY (PRINT NAME): Daniel Kople

SIGNATURE: 2/10/16

PAGE ____ OF ____

-

| CLIENT: | | | | | WELL | ID: | MW | -11 | _ | |
|---------------------------|-----------|--------|---|-----|------|--------|-------|------|--------------|--|
| PROJECT #: (10psey | | | | | ADDR | ESS:_2 | 002 - | 2004 | (represe Are | |
| DATE: 2/10/16 | | | | | SAMP | LING C | REW: | Dan | + Ashlyn | |
| DEPTH TO WATER (DTW) | 18 | .10 ft | _ | D | TB=5 | 50.194 | 4 | | / | |
| CASING DIAMETER (inches): | <u>×2</u> | 3 | 4 | 4.5 | 5 | _6 | _8 | _12 | Other | |

PRODUCT PRESENT:_____ PRODUCT THICKNESS:_____

| TIME (2400 Hr) | VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE °F/°C | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
|-------------------|------------------|---------------|-------------------------|----------------------|---------------|-----|--------------------|-------------------|
| 1250 | 58 | 6,77 | 0.888 | 19.05°C | 8.10 | 169 | 10.4 | Clear |
| 12.55 | 59.5 | 6.86 | 0,880 | 20.4100 | 2.47 | 118 | 11.4 | Chear |
| 1300 | 60.7 | 6.89 | 0.818 | 19.76°C | 2,40 | 110 | 7.4 | Clean |
| 1305 | 61.8 | 6.91 | 0.882 | 20,16°C | 2,25 | 105 | 3.0 | Chean |
| 1310 | 62.8 | 6.89 | 0.882 | 20.3200 | 2.21 | 102 | 0.0 | Chear |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):____

| | PURGING E | QUIPMENT | | SAMPLING I | EQUIPMENT |
|--------|---------------------|--------------------------|---|---------------------|--------------------------|
| X | 2" Submersible Pump | Bailer (Teflon®) | X | 2" Submersible Pump | Bailer (Teflon®) |
| - 4 | Centrifugal Pump | Bailer (PVC) | | DDL Sampler | Bailer (Stainless Steel) |
| | 4" Submersible Pump | Bailer (Stainless Steel) | - | | Peristaltic Pump |
| | Dedicated | Peristaltic Pump | | Disposable Bailer | Dedicated Bailer |
| Other: | | | | Other: | |

WALL CONDITION: _____ GROUT: _____ MANWAY: _____ LOCKING CAP: _____ LOCK: _____ REMARKS: Horiba - 21780

Monsioon GW Pump -16773

COMPLETED BY (PRINT NAME): Davie Kope c

| SIGNATU | RE: | m | 1 h | |
|---------|-----|---|-----|--|
|---------|-----|---|-----|--|

PAGE _____ OF _____

| CLIENT: | WELL ID: MW-02 |
|---------------------------|----------------------------------|
| PROJECT #: Cropsey | ADDRESS: 2002-2004 (ropsey Ave. |
| DATE: 2/10/16 | SAMPLING CREW: Dun + Achlyn |
| DEPTH TO WATER (DTW) | 19.28 ft DTB= 27.39 ft |
| CASING DIAMETER (inches): | <u>X2 3 4 4.5 5 6 8 12</u> Other |

PRODUCT PRESENT: _____ PRODUCT THICKNESS: _____

| VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE °F/°C | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
|------------------|--------------------------|---|---|--|---|---|--|
| 25 | 6.41 | 0.69 | 19.28°C | 1,46 mg/L | 189 | 32.2 | Clear |
| 26 | 6.52 | 0.675 | 19.70°C | 1,99 | 160 | 25,2 | Clear |
| 27 | 6,51 | 0,667 | 20.60°C | 1.96 | 152 | 20.2 | Clear |
| 28 | 6.54 | 0.672 | 20.70°C | 2,02 | 148 | 21.4 | Chear |
| | (gal.) 25 26 27 | $\begin{array}{c c} (gal.) & (units) \\ \hline 2.5 & 6.41 \\ \hline 2.6 & 6.52 \\ \hline 2.7 & 6.51 \\ \hline 0.0 & 6.54 \\ \hline \end{array}$ | (gal.) (units) (ms/cm) 2.5 6.41 0.691 2.6 6.52 0.675 2.7 6.51 0.667 | (gal.) (units) (ms/cm) °F/°C 2.5 6.41 0.691 19.28 °C 2.6 6.52 0.675 19.70 °C 2.7 6.51 0.667 20.60 °C | (gal.) (units) (ms/cm) °F/°C (ppm) 2.5 6.41 0.691 19.28°C 1.46 mg/L 2.6 6.52 0.675 19.70°C 1.99 2.7 6.51 0.667 20.60°C 1.96 | (gal.) (units) (ms/cm) °F/°C (ppm) 2.5 6.41 0.69 19.28 °C 1.46 mg/L 189 2.6 6.52 0.675 19.70 °C 1.99 160 2.7 6.51 0.667 20.60 °C 1.96 152 | (gal.)(units)(ms/cm) $\circ T/\circ C$ (ppm)(NTU)2.56.41 0.691 $19.28°C$ $1.46 m_3/L$ 189 32.2 2.66.52 0.675 $19.70°C$ 1.49 160 25.2 2.76.51 0.667 $20.60°C$ 1.96 152 20.2 |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):_

| | PURGING E | QUIPMENT | | SAMPLING I | EQUIPMENT |
|--------|---------------------|--------------------------|---|---------------------|--------------------------|
| X | 2" Submersible Pump | Bailer (Teflon®) | X | 2" Submersible Pump | Bailer (Teflon®) |
| | Centrifugal Pump | Bailer (PVC) | _ | DDL Sampler | Bailer (Stainless Steel) |
| | 4" Submersible Pump | Bailer (Stainless Steel) | | | Peristaltic Pump |
| | Dedicated | Peristaltic Pump | | Disposable Bailer | Dedicated Bailer |
| Other: | | | | Other: | |

WALL CONDITION: _____ GROUT: _____ MANWAY: _____ LOCKING CAP: _____ LOCK: _____ REMARKS: Moriba - 21780 Monsoon 6W Pump - 16773 .

COMPLETED BY (PRINT NAME): Daniel Kopec

| SIGNATURE: | mont he |
|------------|---------|
| DATE: 2 | 110/16 |

PAGE ____ OF ____

and the second

| CLIENT: | - | | | WEL | LID: MW | -3 | | |
|-------------------|------------------|----------------|-------------------------|----------------------|---------------|----------|--------------------|-------------------|
| PROJECT #: | Cropser | | | ADD | RESS: | | | |
| DATE: 2/ | DATE: 2/10/16 | | | | PLING CREV | V: | | |
| DEPTH TO | WATER (DTW | D _1° | 1.64 | 27.93 | オート | 5 314 = | :15 | |
| CASING DL | AMETER (incl | nes): <u>2</u> | 3 4 | 4.5 _5 | 6 | 8 12 | Other | |
| | | | | Stort | purge | 1545 | | |
| | | | | limite | d drim | space, p | urged 8 gi | Mons. |
| PRODUCT PRE | SENT: | _ PRODUCT T | HICKNESS: | | | | <u>.</u> | |
| TIME (2400 Hr) | VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE °F/°C | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
| 1555 | 8 | 6.82 | 0.500 | 14.89 | 11.30 | 234 | JQ.0 | clear |
| 1600 | 8.5 | 6.50 | 0.490 | 16.11 | 12-21 | 255 | 3.5 | |
| 1605 | 9.0 | 6.44 | 0.480 | 16.55 | 12.10 | 246 | 45.7 | |
| 1610 | 9.5 | 6.47 | 0,483 | 16.64 | 11.88 | 251 | 10.1 | |
| 1613 - | - Sam | ple. col | lected - | _ | 1. | | | |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):_

| PURGING I | QUIPMENT | SAMPLING | EQUIPMENT |
|---------------------|--------------------------|---------------------|--------------------------|
| 2" Submersible Pump | Bailer (Teflon®) | 2" Submersible Pump | Bailer (Teflon®) |
| Centrifugal Pump | Bailer (PVC) | DDL Sampler | Bailer (Stainless Steel) |
| 4" Submersible Pump | Bailer (Stainless Steel) | | Peristaltic Pump |
| Dedicated | Peristaltic Pump | Disposable Bailer | Dedicated Bailer |
| ther: | | Other: | |

| WALL CONI | | GROUT: | MANWAY: | LOCKING CAP: | LOCK: | |
|-----------|---------------------------------------|--------|---------|--------------|-------|--|
| REMARKS: | Honba | 021674 | | | | |
| | Monsoon | 110660 | | | | |
| | A A A A A A A A A A A A A A A A A A A | | | | | |
| | | | | | | |

COMPLETED BY (PRINT NAME):____

SIGNATURE:_____

DATE:

PAGE ____ OF ____

| | OROCILE | WELL ID: MW 45 |
|---|----------------|--|
| CLIENT: WOPCAY PROJECT #: WOPCAY DATE: 2- 10/14 DEPTH TO WATER (DTW) | 18:54 | ADDRESS: |
| CASING DIAMETER (inches): | <u>X 2 3 4</u> | 4.5 _5 _0 _0 _0 1237 Stanto pringel xt=5.6 34xt=16.7 |

PRODUCT THICKNESS: PRODUCT PRESENT:_____

COLOR TURBIDITY ORP TEMPERATURE D.O. (visual) CONDUCTIVITY (NTU) pH VOLUME °F(°G (ppm) TIME (ms/cm) (units) (gal.) (2400 Hr) 01.1 clear 21.9 245 13.71 1.65 14 15 01 clear 1255 216 11 . 15.46 1.(0) 1300 6.40 16 279 0.6 14.96 13.01 6.22 18.5 61 1305 0.0 10.15 274 16-27 1.59 19 5 6.17 1310 10.95 0.0 23 14 25 1 4.6 0 20.00 1315

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):_

| | PURGING EC | QUIPMENT | SAMPLING I | |
|-------|---------------------|--------------------------|---------------------|--------------------------|
| I | | Bailer (Teflon®) | 2" Submersible Pump | Bailer (Teflon®) |
| K | 2" Submersible Pump | Bailer (PVC) | DDL Sampler | Bailer (Stainless Steel) |
| | Centrifugal Pump | Bailer (Stainless Steel) | | Peristaltic Pump |
| | 4" Submersible Pump | Peristaltic Pump | Disposable Bailer | Dedicated Bailer |
| Other | | | Other: | |

LOCKING CAP: LOCK: MANWAY:____ GROUT: WALL CONDITION: 295 21 5000 HonDa REMARKS: 022611 Monsoon COMPLETED BY (PRINT NAME): ASTULIA an SIGNATURE:

110716 DATE:_ 2

| Tim | ¥ | pt | Cond. | Tump. | 70 | A. a. | | |
|------|----|----------------|-------|-------|-------|-------|------|-------------|
| 1320 | 1 | (e.6) (e-30 | 1.44 | 11.8 | 8.73 | 247 | TUM. | Color |
| 1330 | 22 | 6.19 | 1.60 | 13:79 | 13.92 | | 6.0 | |
| 335 | 23 | 4.65 | 1.59 | 14.59 | | 294 | | er Tichelar |

- collected sample @ 1335 -

| CLIENT: WOPRE | WELL ID: MW - 4I |
|--------------------------------------|---------------------------|
| PROJECT #: | ADDRESS: |
| DATE: 2/10/14 | SAMPLING CREW: _ AW |
| DEPTH TO WATER (DTW) | 37.45 7=37.3 |
| CASING DIAMETER (inches): $2 3 (14)$ | <u>4.5 5 6 8 12</u> Other |

PRODUCT PRESENT:_____ PRODUCT THICKNESS:_____

| TIME (2400 Hr) | VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE °F/°C | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
|-------------------|------------------|---------------|-------------------------|----------------------|---------------|-----------|--------------------|-------------------|
| 1916 | 21 | 6.77 | 1.16 | 8.22 | 14.40 | 12:00 128 | 2001 | Clear |
| 421 | 22 | 6.74 | 0.600 | 8.40 | 14.48 | JANK139 | 237 | |
| 1426 | 22.5 | 6.79 | 1.63 | 8.39 | 14.95 | 258 | 123 | |
| 1431 | Actor Jas | 6.91 | 1.40 | 14.57 | 12.49 | 284 | 47.7 | |
| 14 34 | 24 | ce. 40 | 1.37 | 15.70 | 12.00 | 186 | 27.9 | 1 |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):____

| PURGING EC | QUIPMENT | SAMPLING I | EQUIPMENT |
|---------------------|----------------------------|---------------------|--------------------------|
| 2" Submersible Pump | Bailer (Teflon®) | 2" Submersible Pump | Bailer (Teflon®) |
| Centrifugal Pump | Bailer (PVC) | DDL Sampler | Bailer (Stainless Steel) |
| 4" Submersible Pump | Bailer (Stainless Steel) | | Peristaltic Pump |
| Dedicated | Dedicated Peristaltic Pump | | Dedicated Bailer |
| | | Other: | |

MANWAY:_____ LOCKING CAP:_____ LOCK:_____ GROUT:_ WALL CONDITION: 021674 REMARKS: HON DA Mongoon 022611

COMPLETED BY (PRINT NAME):_____

ŧ.

SIGNATURE:_

DATE:_

PAGE ____ OF ____

| | | | Cond. | Temp. | DO | OR | TUND. | Color |
|------|------|-----------|-------|-------|----|----|-------|-------|
| 1441 | 75 | 4.43 | 1.35 | | | | 12.3 | |
| | - 50 | angeli (e | 1445 | | | | | |

| CLIENT: CNOPSUY | WELL ID: MW-5 |
|-----------------------------------|----------------------------|
| PROJECT #: | ADDRESS: 2007 Cropley tre. |
| DATE: 3/18/16 | SAMPLING CREW: 10 |
| DEPTH TO WATER (DTW) | |
| CASING DIAMETER (inches): $4 4$. | <u>5 5 6 8 12</u> Other |
| Start purge @ 10 20 | а 1 |

PRODUCT PRESENT:_____ PRODUCT THICKNESS:_____

| TIME (2400 Hr) | VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE °F°C | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
|-------------------|------------------|---------------|-------------------------|---------------------|---------------|-----|--------------------|-------------------|
| 1037 | 6 | 6.52 | 0.551 | 16.12 | 10.32 | 253 | 58.1 | clear |
| 1042 | 7 | 6.47 | 0-516 | 14.99 | 9.29 | 240 | 39 | clow |
| 1047 | 8 | 6.47 | 0.509 | 17.07 | 9.79 | 235 | 21.7 | char |
| 1052 | 9 | 6.49 | 0.510 | 16.45 | 9.48 | 231 | 12.6 | char |
| 1057 | 10 | 4.51 | 0.528 | 14.98 | 9.97 | 230 | 7.4 | clear |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):__

| | PURGING EQ | QUIPMENT | SAMPLING EQUIPMENT | | | | |
|------------------|--------------------------------------|--------------------------|---------------------|--------------------------|--|--|--|
| 2" Submersible F | 2" Submersible Pump Bailer (Teflon®) | | 2" Submersible Pump | Bailer (Teflon®) | | | |
| Centrifugal Pum | , | Bailer (PVC) | DDL Sampler | Bailer (Stainless Steel) | | | |
| 4" Submersible I | ump | Bailer (Stainless Steel) | | Peristaltic Pump | | | |
| Dedicated | | Peristaltic Pump | Disposable Bailer | Dedicated Bailer | | | |
| her: | | | Other: | | | | |

MANWAY:_____ LOCKING CAP:____ LOCK:__ WALL CONDITION: GROUT: REMARKS: MONSOON "Regular #1" Flow controlle 15 Horiba 355 or hun COMPLETED BY (PRINT NAME): SIGNATURE DATE:

PAGE OF

| Time | vol | pA | Cond. | Temp | 20 | ORP | Turb. | Color | |
|------|------|-----------|-------|-------|--------|-----|-------|-------|--|
| 1102 | 10 | 6.45 | 0.500 | 16.96 | 0.07 | 230 | 6.2 | clian | |
| 1107 | n | 6.48 | 0.498 | 16:74 | 10.41 | 730 | 5.3 | cliar | |
| 1102 | 12 | 4.55 | 0.499 | 16.93 | 9.55 | 227 | 3.7 | clear | |
| 1117 | 12.5 | 4.55 | 0-494 | 17.15 | 9.49 | 227 | 3.2 | clear | |
| | _ | collected | sampl | 13 @ | 1120 - | | | | |

| CLIENT: Cropsey | WELL ID: MW-7 ADDRESS: 2002-2004 (repsey Ave SAMPLING CREW: Dan | | | | | | | | |
|---------------------------|---|----|-----|------|-------|------|------|-------|--|
| PROJECT #: Cropsey | | | | | | _ | | | |
| DATE: 3/18/16 | | 0 | | SAMP | LINGC | REW: | ISMA | | |
| DEPTH TO WATER (DTW) | 20.14 | 1+ | | | | | | | |
| CASING DIAMETER (inches): | χ_2 3 | 4 | 4.5 | 5 | _6 | 8 | 12 | Other | |

PRODUCT PRESENT:___

PRODUCT THICKNESS:

Donnel Kaper

| TIME (2400 Hr) | VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE °F/°C | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
|-------------------|------------------|---------------|-------------------------|----------------------|---------------|--------|--------------------|-------------------|
| 0954 | 6 | 6,20 | 0779 | 17.89 | 8.32 | -169 | 127 | alear |
| 0959 | 7.5 | 6.27 | 0.750 | 19.05 | - 8,48 | 101 | 138 | Clear |
| 1004 | 9 | 134 | 0.742 | 18.91 | 5,40 | 10 102 | 32,3 | Clear |
| 1009 | 10 | 6.39 | 0.731 | 18,40 | 4.83 | 93 | 17.8 | Cheur |
| 1014 | 11 | 6.38 | 0.734 | 18.45 | 8.24 | 88 | 15.2 | Clear |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):___

| | PURGING E | QUIPMENT | SAMPLING I | EQUIPMENT | | |
|-------|--------------------------------------|--------------------------|---------------------|--------------------------|--|--|
| V | 2" Submersible Pump Bailer (Teflon®) | | 2" Submersible Pump | Bailer (Teflon®) | | |
| A | Centrifugal Pump | Bailer (PVC) | DDL Sampler | Bailer (Stainless Steel) | | |
| | 4" Submersible Pump | Bailer (Stainless Steel) | | Peristaltic Pump | | |
| | Dedicated | Peristaltic Pump | Disposable Bailer | Dedicated Bailer | | |
| Other | Louisites | | Other: | | | |

LOCK: LOCKING CAP: MANWAY: GROUT: WALL CONDITION: 3 599 10 yme 1945. We Pu-upo REMARKS: Dury ted ask 518 30

COMPLETED BY (PRINT NAME):____

SIGNATURE:_ DATE:

PAGE ____ OF ____

| me | VOL | Ph | Cond | Temp | | ORP | | |
|-----|-------|--|---|--|---|-----|------|-------------------------|
| 019 | 12 | 6.34 | .0.739 | 19.02 | 7.38 | 89 | 10,5 | 0 |
| | | | | 1 | * | | | |
| | | | | | | | | |
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R.

| CLIENT: Crepsey | WELL ID:6 |
|--|--------------------------------------|
| PROJECT #: (napley | ADDRESS: 2002-2004 (upper Ave BK, NY |
| DATE: 3/18/16 | SAMPLING CREW: 10 dm |
| DEPTH TO WATER (DTW) 20.05 | Et |
| CASING DIAMETER (inches): $\underline{\times 2}$ 3 | <u>4 4.5 5 6 8 12</u> Other |

PRODUCT PRESENT: _____ PRODUCT THICKNESS: _____

| TIME (2400 Hr) | VOLUME (gal.) | pH (units) | CONDUCTIVITY (ms/cm) | TEMPERATURE | D.O. (ppm) | ORP | TURBIDITY (NTU) | COLOR (visual) |
|-------------------|------------------|---------------|-------------------------|-------------|---------------|-----|--------------------|-------------------|
| 1100 | 6 | 6.78 | 0,656 | 17.90 | 5,44 | 178 | 106 | Musky |
| 1105 | 6 | 6.59 | 0.654 | 18,36 | 6.99 | 185 | 53.5 | Clear |
| 1110 | 7 | 6.65 | 0.655 | 17.41 | 7.90 | 190 | 25,4 | aleur |
| 1115 | 8 | 6.52 | 0.646 | 18.23 | 4.07 | 187 | 16.4 | Clear |
| 1120 | 9 | 6.50 | 0.646 | 18.31 | 7,14 | 180 | 11.1 | Clear |

FIELD QC SAMPLES COLLECTED AT THIS WELL (i.e, FB-1, XDUP-1):___

| | PURGING E | QUIPMENT | SAMPLING EQUIPMENT | | | | |
|--------|---------------------|--------------------------|--------------------|---------------------|--------------------------|--|--|
| X | 2" Submersible Pump | Bailer (Teflon®) | | 2" Submersible Pump | Bailer (Teflon®) | | |
| | Centrifugal Pump | Bailer (PVC) | | DDL Sampler | Bailer (Stainless Steel) | | |
| | 4" Submersible Pump | Bailer (Stainless Steel) | | | Peristaltic Pump | | |
| | Dedicated | Peristaltic Pump | | Disposable Bailer | Dedicated Bailer | | |
| Other: | | | | Other: | | | |

WALL CONDITION: GROUT: MANWAY: LOCKING CAP: LOCK: 4 gu urue 62 unsed REMARKS: 3x volume n Jule-en Que 1 mm. COMPLETED BY (PRINT NAME):_ unte aper SIGNATURE:_ DATE:_

PAGE ____ OF ____

| Time | Volume | pH 1 | Lond | Temp | DO 1 | ORP | Turlo | Color |
|------|--------|------|--|--|------|-----|-------|-------|
| 1125 | 10 | 6.58 | 0.644 | 18.10 | 3,18 | 182 | 8.7 | Chear |
| 1130 | 11-1 | 6,56 | 0.646 | 18.3) | 3.09 | 184 | 6.6 | Clear |
| 1135 | 12 | 6.56 | 0.645 | 18.37 | 2.28 | 189 | 5.2 | Clear |
| 1140 | 13 | 6.55 | 0.649 | 17.88 | 4,39 | 193 | 4.3 | Clev- |
| 1145 | | | | | | 1 | | |
| 7 | Λ Λ | | and other second se | and the second | | | | |
| | | | | | |) | | I |



| Client: | AMA Cro | psey | | Location (Si | te/Facility Nam | ie): | (vopse) | , Are, | Brookly | 5 |
|-----------------------------|---|--------------------|-------------|-------------------|------------------|-------------|-------------|---|-------------|---------------------------|
| | |) F | ield Perso | nnel: Da | n Kople | | | Weather: | Sunny | High 80 |
| ate: | | | | | | | | | / | |
| Vell Number | c . | MW-5I | | Evacuation | Method: | | | | | |
| Sampling Me | ethod: | Low Flo | h- | Purging Dev | vice (pump type | e): _ | 5.5. | Monsoo | 5 | |
| Vell Informa | tion: | | Screen Ber | nin | | ft. | *Measurem | ents Taken | From | |
| Depth of We | * | | | | | ft. ft. | | | Top of Wel | l Casing ective Casing |
| Depth of Wa | tor * | 19.64 | Screen Lt | 10 | | π. | 1. j | | (Other, Spe | |
| | | - A. | - | | 45 | 4 | | Well Casin | g size _ | 211 |
| _ength of W | ater Column | | Pump Set | @ Mid-Point | | n | | Wen odoni | 9 0120 _ | |
| Water | | | | | | | | | | |
| parameters The last 3 re | adings must sta | bilize to criteria | a below. Sa | ampling rate | of 0.2-0.25 lite | ers/minute. | | | | |
| Elapsed time | es will reflect the | e actual exact t | me of read | lings. | | 10% | 0.3 mg/l | 200-500 | N/A | N/a |
| Criteria | 0.3 ft | 3% | 0.1 | 10mv Oxidation | 3% | 10% | Dissolved | 200-500 | | |
| Elapsed | Dept To Water | Temperature | | Reduction | Conductivity | Turbidity | Oxygen | Flow Rate | | Odor |
| Time | ft | degree C | pН | m. volts | U ohm/cm | (NTU) | (mg/l) | (ml/min) | T T | Apparent |
| 1247 | | 19.61 | 6.07 | 163 | 0.913 | 0.0 | 7.32 | 280 | Clear | None |
| 1252 | | 18.52 | 6.23 | 152 | 0.924 | 0.0 | 5.23 | 480 | Cleur | None |
| 1257 | | 18.40 | 6.38 | 146 | 0.928 | 0.0 | 4.43 | 220 | Clear | None |
| 1302 | | 18,12 | 6.42 | 143 | 0.933 | 0.0 | 4.07 | 480 | Clear | some |
| 1307 | | 18.23 | 6.45 | 143 | 0.933 | 0.0 | 3.84 | 440 | Cleur | None |
| 1312 | | 18.43 | 6.45 | 142 | 0.935 | 0.0 | 3,52 | 320 | Clear | None |
| 1317 | | 18.20 | 6.45 | 143 | 0.936 | 0,0 | 3.43 | 480 | Chevr | None |
| 1319 | - saw | ple- | | | | | | | | |
| | | 1 | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | | | | | | | | |
| | | | O / CALLE | | | | | | | |
| | ple: Mw-5 | I , MW - 10 | o(Field 1 | Jup.) | | | | | | |
| | t: <u>1117</u> | | - | | Total volume | of purged | water remo | ved: | | |
| Time Colle | and the second se | | _ | | Total volume | Physical | appearance | e at samplir | ng | |
| Physical ap | opearance at sta | Clear | | | | i nyorodi | Color | Cheur | - | |
| | Color | | | | | | Odor | None | | |
| Sheen/Fre | Odor | Vone | _ | | | Sheen/F | ree Product | the second se | | |
| NOTES: | | pural @ | 1117 | | | Sam | phe to ku | n Q 13 | 19 | |
| NOTES: | Stort | Dune @ | 1247 | | | | | | | |
| | Star | T. O M | w @ 12 | 47 | | | | | | |
| 1 | | IN THE L | 0 | 1 | | | | | | |



| Client: | CATOODEM | | | Location (Si | te/Facility Nam | e): | | | | |
|-----------------|--------------------|-------------------------|-------------------------|------------------------|------------------|------------|-----------------------|--------------|---------------------------------------|-----------------|
| | 2 10 hr | | | | | | | Weather: | ~ 90°F | forny. |
| Date: | 7/8/16 | | | nnel: AN | | | 10 | - | , | 0 |
| Well Number: | | MW-8 | | Evacuation | Method: | sub. | | | | |
| Sampling Meth | nod: | 10W-Plou |) | Purging Dev | vice (pump type | e): _ | Monsoe | M | | |
| Well Informatio | on: | 0.00 | Screen Beg | gin 16. | | | *Measurem | ents Taken | | " O size |
| Depth of Well* | | | Screen Eng Screen Lt | | 8 | ft. ft. | | | Top of We Top of Pro (Other, Sp | otective Casing |
| Depth of Wate | er * | 18.42 | | | | | l | | (Other, Sp | |
| Length of Wat | ter Column | 8.38 | Pump Set (| @ Mid-Point | | ft. | | Well Casin | g size | 2" |
| Water | | | | | | | | | | |
| parameters: | dinge must sta | abilize to criteria | below. Sa | mpling rate | of 0.2-0.25 lite | rs/minute. | | | | |
| Elapsed times | s will reflect the | e actual exact t | me of read | ings. | | | | | | N/- |
| Criteria | 0.3 ft | 3% | 0.1 | 10mv | 3% | 10% | 0.3 mg/l Dissolved | 200-500 | N/A | N/a |
| | Dept | Temporatura | | Oxidation Reduction | Conductivity | Turbidity | Oxygen | Flow Rate | Color | Odor |
| Elapsed Time | To Water ft | Temperature degree C | pН | m. volts | U ohm/cm | (NTU) | (mg/l) | (ml/min) | Apparent | |
| 1040 | 18.31 | 16.81 | 6.27 | -56 | 1.33 | 0.0 | 0.00 | 488 | clear | None |
| 1045 | 18.31 | 16.72 | 6.26 | -58 | 1.33 | 0.0 | 0.00 | 480 | | |
| 1050 | 18.31 | 16.92 | 6.26 | -60 | 1.32 | 0.0 | 0.00 | 480 | | |
| 1055 | 18.31 | 17.12 | 6.78 | -42 | 1.32 | 0.0 | 0.00 | 420 | | |
| 001100 - | -san | iple - | | | | | | | | |
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| | | | | | | | | | 1 | 1 |
| Water Samp | | | | | | | | | | |
| Pump Start: | | | | | Total volume | of purged | water remo | ved: | | |
| Time Collect | | | - | | rotar volume | Physical | appearance | e at samplin | g | |
| Physical app | earance at sta | hrown | | | | , njelour | Color | | | |
| | Color Odor | Nerry. | - | | | | Odor | | | |
| Sheen/Free | | TUDY M. | - | | | Sheen/F | ree Product | | | |
| NOTES: | | | | | | | 0 | C a A | 20 | |
| | MS | MSD COM | rected | | | | wrad | ~ 5.0 9 | sel. | |
| | | soon pum | P # R | 10931 | | | | | | |
| | itor | 10a 2119 | 1 | | | | | | | |



| Client: | AMA | | | Location (Si | te/Facility Nam | ie): | Cropse | y | | |
|--------------|---------------------|---------------------|-------------------------|------------------------|------------------|------------|-----------------------|---------------|-------------------------------|----------------|
| | 010/11 | | | A. | | | | Ø Weather: | ~90°F | SOMMU |
| Date: | 718/16 | | ield Perso | nnel: /\/ | V | | | - | | 0 |
| Vell Numbe | r. | MW-9 | | Evacuation | Method: | Sub. | | | | |
| Sampling Me | ethod: | LOW-PLOI | N | Purging Dev | vice (pump type | e): _ | Monso | m | | |
| Well Informa | ation: | | Screen Beg | gin 🚳 [| 6.77 | ft. | *Measurem | ents Taken | From | |
| Depth of We |) * | 26.77 | Screen Eng Screen Lt | 1 26.7° | | ft. ft. | | | and the product of the second | tective Casing |
| Depth of Wa | ater * | 18.27 | | | | | | | (Other, Spe | ecify) |
| Length of W | ater Column | | Pump Set (| @ Mid-Point | | ft. | | Well Casin | g size | |
| Water | | | | | | | | | | |
| parameters | adings must sta | abilize to criteria | a below. Sa | mpling rate | of 0.2-0.25 lite | rs/minute. | | | | |
| Elapsed tim | es will reflect the | e actual exact t | ime of read | ings. | | | 0.0 | 200 500 | N/A | N/a |
| Criteria | 0.3 ft | 3% | 0.1 | 10mv | 3% | 10% | 0.3 mg/l Dissolved | 200-500 | IN/A | 14/0 |
| Elapsed | Dept To Water | Temperature | | Oxidation Reduction | Conductivity | Turbidity | Oxygen | Flow Rate | Color | Odor |
| Time | ft | degree C | pН | m. volts | U ohm/cm | (NTU) | (mg/l) | (ml/min) | Apparent | Apparent |
| 1220 | 18.13 | 28.14 | 6.34 | -89 | 2.25 | 0.0 | 0.00 | 520 | clean | none |
| 1225 | 18-10 | 27.81 | 6.38 | -94 | 5-38 | 0.0 | 0.00 | 260 | | 1 |
| 1230 | 18.10 | 20.79 | 6-38 | -96 | 2.29 | 1.5 | 0.00 | 200 | | |
| 1235 | 18.10 | 27.73 | 6.39 | -99 | 2.27 | 0.1 | 0.0 | 280 | | 1 |
| 1240 | | 0 | 4.01 | | 0407 | | | | | |
| 1940 | - sam | pre | | | | | | | | |
| | - | | | | | | | | | |
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| Mada Cam | | | | | | | 10 | | | |
| Water San | | | | | | | | | | |
| Pump Star | | 0 | -0 | | Total volume | of purged | water remov | ved: ~ ° | 7 gal | |
| Time Colle | | | - | | | Physical | appearance | e at samplir | | |
| Physical a | ppearance at sta | ONRIA | | | | | Color | clear | | |
| | Color | 1000 | _ | | | | Odor | none | | |
| Sheen/Fre | Odor a Broduct | none | - | | | Sheen/F | ree Product | | | |
| | | on pump | # RIG | 1931 | | | | | | |
| NOTES: | Mons | | 7 101 | | | D | urged ~ | 6.0 900 | | |
| | forib | a and m | trover | | | 1 | ha | 4 | | |
| | 11-1 | N-one con | VIA GION | | | | 191 | | | |
| 4 | | | | | | | | | | |



| lient: | AMA (| ropsey | l | ocation (Sit | e/Facility Nam | e): | Crop | | Brook | / |
|-----------------------------|----------------|--|--------------|------------------------|--------------------|------------|---|--|---------------|----------|
| _ | -10/11 | 1 / | ald Dame | | n Kopec | | 1 | Neather: | Sunnu | 805 |
| te: _ | 118/16 | F | ield Persol | | n reppeo | | | 100-100 | | |
| ell Number: | | MW-10 | | Evacuation I | Method: | | | | | |
| ampling Meth | nod: | Low Flow | / | Purging Dev | rice (pump type | e): | 5.5 M | lonsoon | | |
| /ell Informatio | on: | | Paragan Para | in | | ft. | *Measurem | ents Taken | From | |
| epth of Well* | ÷ | 2700 0 | Creen Beg | 1 | | ft. | F | | Top of Wel | Casing |
| | | 10 110 | Screen Lt | 10 | | ft. | | | Other, Spe | |
| epth of Wate | er* | 19.46 | | | | | | | | 2" |
| ength of Wat | ter Column | F | Pump Set (| Mid-Point | 22 | ft. | | Well Casing | size _ | 2 |
| /ater arameters: | | | | | | | | | | |
| he last 3 rea | dings must sta | abilize to criteria | below. Sa | mpling rate | of 0.2-0.25 lite | rs/minute. | | | | |
| Criteria | 0.3 ft | e actual exact til | 0.1 | 10mv | 3% | 10% | 0.3 mg/l | 200-500 | N/A | N/a |
| | Dept | Temperature | | Oxidation Reduction | Conductivity | Turbidity | Dissolved Oxygen | Flow Rate | Color | Odor |
| Elapsed Time | To Water ft | degree C | рН | m. volts | U ohm/cm | (NTU) | (mg/l) | and the second sec | Apparent | Apparent |
| 1015 | | 17,01 | 6.13 | 163 | 0.919 | .1.1 | 2.39 | 240 | Cleor | None |
| 1020 | | 17.25 | 6.08 | 163 | 0.924 | 1,5 | 2.27 | 220 | Clear | None |
| 1025 | | 17.42 | 6.03 | 164 | 0.926 | 1.7 | 2.05 | 300 | C-leo- | Nore |
| 1020 | | 16.85 | 5.98 | 168 | 0.930 | 0.5 | 2.16 | 400 | Clea- | None |
| 1032 | 50 | mple- | | | | | | | | |
| 1034 | 5 | of of a | | | | | | | | |
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| | | | | | | | | | | |
| .4. | | | | | | | | | | |
| Pump Start: Time Collect | | - |]~ກ | | Total volume | Physical | water remo appearanc Color Odor Free Produc | e at samplir <u>Claur</u> <u>None</u> | a) ng | |
| Sheen/Free | | | 0001 | | C | | | | | |
| NOTES: | Start Stop | purge (2) prove (2) Jon-flow (2) | 1015 - | | Sample ip to Ho | | 0 1032 | | | |
| 5. | Marila | nsoon 2" Pu → 35330 → 20532 | | 134 | Som | e equi | pmerit i | Por Min | - 5Ī - 100 | |

Appendix G

Groundwater Data Analytical Reports



Section 4

Sample Results

Report of Analysis





| | | | - 1 | | J | | 6 |
|--|------------------------------|------------------------------|--------------------------|-----------------|----------------------|--|-----------------------------------|
| Client Sar Lab Samp Matrix: Method: Project: | e ID: JC14 AQ EPA | 4061-1 - Ground Wa 624 | ater sey Avenue, Bro | ooklyn, N | I F | Date Sampled: (Date Received: (Percent Solids:) | |
| Run #1 ^a Run #2 | File ID U203278.D | DF 1 | Analyzed 02/11/16 | By NH | Prep Date n/a | Prep Batch n/a | Analytical Batch VU9349 |
| Run #1 | Purge Volun 5.0 ml | ne | | | | | |

Report of Analysis

Run #2

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | 0.89 | 1.0 | 0.091 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 0.72 | 1.0 | 0.14 | ug/l | J |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

11 of 963

ACCUTEST JC14061

SGS

E = Indicates value exceeds calibration range

| Client Samp Lab Sample Matrix: Method: Project: | | venue, Broo | klyn, NY | | Date | Sampled: Received: ent Solids: | 02/10/16 02/10/16 n/a |
|---|-----------------------------|-------------|----------|------|-------|--------------------------------------|-----------------------------|
| VOA PPL I | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.12 | ug/l | | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylenes (total) | ND | 1.0 | 0.22 | ug/l | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 (SUR) | 109% | | 72-1 | 25% | | |
| 2037-26-5 | Toluene-D8 (SUR) | 102% | | 78-1 | 19% | | |
| 460-00-4 | 4-Bromofluorobenzene (SUR) | 104% | | 74-1 | 15% | | |
| 1868-53-7 | Dibromofluoromethane (S) | 105% | | 79-1 | 20% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





SGS

| Client Sar Lab Samj Matrix: Method: Project: | ole ID: JC140 AQ - EPA | Ground Wa 624 | iter sey Avenue, Bro | ooklyn, NY | Da Pe | te Sampled: 0 te Received: 0 rcent Solids: r | |
|--|------------------------------|------------------|-------------------------|------------|-----------|--|----------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| | rne ID | | | | | | i indi y ticui Dutti |
| Run #1 ^a | U203279.D | 1 | 02/11/16 | NH | n/a | n/a | VU9349 |
| Run #1 ^a Run #2 ^a | | 1 10 | • | v | - | - | • |

Report of Analysis

Run #1 5.0 ml

Run #2 5.0 ml

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|------------------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | 0.49 | 1.0 | 0.091 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | 0.72 | 1.0 | 0.19 | ug/l | J |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 6.7 | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 740 ^b | 10 | 1.4 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





Page 1 of 2

13 of 963

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

| Client Samp Lab Sample Matrix: Method: Project: | | zenue, Brook | lyn, NY | | Date | Sampled: Received: ent Solids: | 02/10/16 02/10/16 n/a |
|---|--|-----------------------|--------------------------|------------------------------|------------------------------|--------------------------------------|-----------------------------|
| VOA PPL I | ist | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 75-69-4 75-01-4 1330-20-7 | Trichloroethene Trichlorofluoromethane Vinyl chloride Xylenes (total) | 9.7 ND ND ND | 1.0 2.0 1.0 1.0 | 0.12 0.20 0.13 0.22 | ug/l ug/l ug/l ug/l | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | | |
| 17060-07-0 2037-26-5 460-00-4 | 1,2-Dichloroethane-D4 (SUR) Toluene-D8 (SUR) 4-Bromofluorobenzene (SUR) | 100% | 109% 102% 102% | 72-1 78-1 74-1 | 19% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

104%

105%

79-120%

(b) Result is from Run# 2

1868-53-7

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



14 of 963

ACCUTEST JC14061

Page 2 of 2

4.2 **4**

| Lab Samı Matrix: Method: Project: | AQ - EPA e | JC14061-3 AQ - Ground Water EPA 624 2002-2024 Cropsey Avenue, Brooklyn, NY | | | | ate Sampled:02ate Received:02orcent Solids:n/ | | |
|--|------------------------|---|----------------------|-----|------------|---|------------------|--|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| | | | 00/11/16 | NH | _ | , - | 1110240 | |
| Run #1 ^a | U203280.D | 1 | 02/11/16 | INT | n/a | n/a | VU9349 | |
| Run #1 ^a Run #2 ^a | U203280.D U203293.D | 1 10 | 02/11/16 02/12/16 | NH | n/a n/a | n/a n/a | VU9349 VU9350 | |

Report of Analysis

Run #1 5.0 ml

Run #2 5.0 ml

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|------------------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | 0.62 | 1.0 | 0.091 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 1.5 | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 249 ^b | 10 | 1.4 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

15 of 963

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

| Client Sample ID:MW-2Lab Sample ID:JC14061-3Matrix:AQ - Ground WaterMethod:EPA 624Project:2002-2024 Cropsey Ave | | | klyn, NY | | Date | Sampled: Received: ent Solids: | 02/10/16 02/10/16 n/a |
|---|--|-----------------------|----------------------------|------------------------------|------------------------------|--------------------------------------|-----------------------------|
| VOA PPL I | ist | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 75-69-4 75-01-4 1330-20-7 | Trichloroethene Trichlorofluoromethane Vinyl chloride Xylenes (total) | 2.4 ND ND ND | $1.0 \\ 2.0 \\ 1.0 \\ 1.0$ | 0.12 0.20 0.13 0.22 | ug/l ug/l ug/l ug/l | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 2037-26-5 460-00-4 | 1,2-Dichloroethane-D4 (S Toluene-D8 (SUR) 4-Bromofluorobenzene (S | 102% | 108% 100% 106% | 78-1 | 25% 19% 15% | | |

Report of Analysis

(a) (pH=5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

104%

104%

79-120%

(b) Result is from Run# 2

1868-53-7

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



16 of 963



| | | | - L | | J | | |
|-----------|----------------|------------|-----------------|------------|-----------|------------------|------------------|
| Client Sa | mple ID: MW-2 | | | | | | |
| Lab Sam | ple ID: JC140 | 61-3 | | | Da | ate Sampled: 02 | 2/10/16 |
| Matrix: | - AQ - (| Ground Wa | ater | | Da | ate Received: 02 | 2/10/16 |
| Method: | EPA 6 | 25 EPA | 625 | | Pe | rcent Solids: n/ | a |
| Project: | 2002-2 | 2024 Crops | sey Avenue, Bro | ooklyn, NY | Y | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | F154587.D | 1 | 02/12/16 | SD | 02/11/16 | OP91141 | EF6502 |
| Run #2 | | | | | | | |
| | Initial Volume | Final V | olume | | | | |
| Run #1 | 980 ml | 1.0 ml | | | | | |

Report of Analysis

Run #2

ABN PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|-----|------|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 5.1 | 0.97 | ug/l | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 5.1 | 1.2 | ug/l | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 5.1 | 1.6 | ug/l | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 5.1 | 1.7 | ug/l | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 5.1 | 0.91 | ug/l | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 5.1 | 0.74 | ug/l | |
| 88-75-5 | 2-Nitrophenol | ND | 2.0 | 1.8 | ug/l | |
| 100-02-7 | 4-Nitrophenol | ND | 10 | 0.86 | ug/l | |
| 87-86-5 | Pentachlorophenol | ND | 5.1 | 1.9 | ug/l | |
| 108-95-2 | Phenol | ND | 2.0 | 0.51 | ug/l | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 2.0 | 1.3 | ug/l | |
| 83-32-9 | Acenaphthene | ND | 1.0 | 0.36 | ug/l | |
| 208-96-8 | Acenaphthylene | ND | 1.0 | 0.39 | ug/l | |
| 120-12-7 | Anthracene | ND | 1.0 | 0.41 | ug/l | |
| 92-87-5 | Benzidine | ND | 20 | 0.29 | ug/l | |
| 56-55-3 | Benzo(a)anthracene | ND | 1.0 | 0.36 | ug/l | |
| 50-32-8 | Benzo(a)pyrene | ND | 1.0 | 0.37 | ug/l | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 1.0 | 0.60 | ug/l | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 1.0 | 0.43 | ug/l | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 1.0 | 0.43 | ug/l | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 2.0 | 0.31 | ug/l | |
| 85-68-7 | Butyl benzyl phthalate | ND | 2.0 | 0.61 | ug/l | |
| 91-58-7 | 2-Chloronaphthalene | ND | 2.0 | 1.0 | ug/l | |
| 106-47-8 | 4-Chloroaniline | ND | 2.0 | 0.41 | ug/l | |
| 218-01-9 | Chrysene | ND | 1.0 | 0.26 | ug/l | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 2.0 | 0.67 | ug/l | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 2.0 | 0.54 | ug/l | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 2.0 | 0.75 | ug/l | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 2.0 | 0.44 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 2.0 | 0.21 | ug/l | |
| 122-66-7 | 1,2-Diphenylhydrazine | ND | 2.0 | 0.46 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 2.0 | 0.16 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

17 of 963

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Client Sample ID: | MW-2 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC14061-3 | Date Sampled: | 02/10/16 |
| Matrix: | AQ - Ground Water | Date Received: | 02/10/16 |
| Method: | EPA 625 EPA 625 | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| Ū | | | |

ABN PPL List

| 106-46-71,4-DichlorobenzeneND2.00.18ug/l $121-14-2$ 2,4-DinitrotolueneND2.00.88ug/l $606-20-2$ 2,6-DinitrotolueneND2.00.57ug/l $91-94-1$ 3,3'-DichlorobenzidineND2.01.2ug/l $53-70-3$ Dibenzo(a, h)anthraceneND1.00.56ug/l $84-74-2$ Di-n-butyl phthalateND2.00.60ug/l $117-84-0$ Di-n-octyl phthalateND2.00.40ug/l $131-11-3$ Dimethyl phthalateND2.00.40ug/l $117-81-7$ bis(2-Ethylhexyl)phthalateND2.00.67ug/l $206-44-0$ FluorantheneND1.00.25ug/l $118-74-1$ HexachlorobenzeneND1.00.46ug/l $118-74-1$ HexachlorobenzeneND1.00.46ug/l $17-72-1$ HexachlorobenzeneND1.00.42ug/l $193-39-5$ Indeno(1,2,3-cd)pyreneND1.00.31ug/l $193-39-5$ Indeno(1,2,3-cd)pyreneND1.00.33ug/l $193-39-5$ NitrobenzeneND2.00.43ug/l $192-03$ NaphthaleneND2.00.44ug/l $192-04$ NaphthaleneND1.00.33ug/l $192-05$ NitrobenzeneND2.00.44ug/l $192-04$ PyreneND1.00.33u | CAS No. | Compound | Result | RL | MDL | Units | Q |
|--|-----------|----------------------------|--------|--------|------|-------|---|
| 606-20-22, 6-DinitrotolueneND2.00.57ug/l91-94-13, 3'-DichlorobenzidineND2.01.2ug/l53-70-3Dibenzo(a, h)anthraceneND1.00.56ug/l84-74-2Di-n-butyl phthalateND2.00.60ug/l117-84-0Di-n-octyl phthalateND2.00.40ug/l117-84-0Di-n-octyl phthalateND2.00.40ug/l131-11-3Dimethyl phthalateND2.00.33ug/l117-81-7bis(2-Ethylhexyl)phthalateND2.00.67ug/l206-44-0FluorantheneND1.00.25ug/l86-73-7FluoreneND1.00.46ug/l118-74-1HexachlorobenzeneND1.00.55ug/l87-68-3HexachlorobutadieneND1.00.42ug/l77-47-4HexachlorobutadieneND5.10.29ug/l193-39-5Indeno(1,2,3-cd)pyreneND1.00.31ug/l98-95-3NitrobenzeneND2.00.43ug/l98-95-3NitrobenzeneND2.00.44ug/l62-75-9n-NitrosodimethylamineND2.00.44ug/l62-75-9n-NitrosodimethylamineND2.00.44ug/l62-75-9n-NitrosodimethylamineND2.00.43ug/l62-75-9n-NitrosodimethylamineND2.00.43ug/l | 106-46-7 | 1,4-Dichlorobenzene | ND | 2.0 | 0.18 | ug/l | |
| 91-94-13,3'-DichlorobenzidineND2.01.2 ug/l 53-70-3Dibenzo(a,h)anthraceneND1.00.56 ug/l 84-74-2Di-n-butyl phthalateND2.00.60 ug/l 117-84-0Di-n-octyl phthalateND2.00.40 ug/l 117-84-0Di-n-octyl phthalateND2.00.33 ug/l 117-81-7bis(2-Ethylhexyl)phthalateND2.00.67 ug/l 206-44-0FluorantheneND1.00.25 ug/l 86-73-7FluoreneND1.00.46 ug/l 118-74-1HexachlorobenzeneND1.00.42 ug/l 87-68-3HexachlorocyclopentadieneND1.00.42 ug/l 67-72-1HexachlorocyclopentadieneND1.00.31 ug/l 91-90-53NitrobenzeneND1.00.33 ug/l 98-95-3NitrobenzeneND1.00.33 ug/l 98-95-3NitrobenzeneND2.00.43 ug/l 62-75-9n-NitrosodinethylamineND2.00.43 ug/l 62-75-9n-NitrosodinethylamineND2.00.48 ug/l 86-30-6N-NitrosodinethylamineND2.00.48 ug/l 86-30-6N-NitrosodiphenylamineND1.00.33 ug/l 129-00-0PyreneND1.00.34 ug/l 120-82-11,2,4-TrichlorobenzeneND1. | 121-14-2 | 2,4-Dinitrotoluene | ND | 2.0 | 0.88 | ug/l | |
| 53-70-3Dibenzo(a,h)anthraceneND1.0 0.56 ug/l 84-74-2Di-n-butyl phthalateND2.0 0.60 ug/l 117-84-0Di-n-octyl phthalateND2.0 0.58 ug/l 84-66-2Diethyl phthalateND2.0 0.40 ug/l 131-11-3Dimethyl phthalateND2.0 0.33 ug/l 117-81-7bis(2-Ethylhexyl)phthalateND2.0 0.67 ug/l 206-44-0FluorantheneND 1.0 0.25 ug/l 86-73-7FluorantheneND 1.0 0.46 ug/l 118-74-1HexachlorobenzeneND 1.0 0.46 ug/l 87-68-3HexachlorobutadieneND 1.0 0.18 ug/l 77-47-4HexachlorocyclopentadieneND 1.0 0.42 ug/l 193-39-5Inden($1,2,3$ -cd)pyreneND 1.0 0.31 ug/l 98-95-3NitrobenzeneND 2.0 0.43 ug/l 98-95-3NitrobenzeneND 2.0 0.43 ug/l 62-75-9n-NitrosodimethylamineND 2.0 0.48 ug/l 86-30-6N-NitrosodiphenylamineND 2.0 0.48 ug/l 129-00-0PyreneND 1.0 0.33 ug/l 129-00-0PyreneND 1.0 0.34 ug/l 120-02-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.35 ug/l 120-03 | 606-20-2 | 2,6-Dinitrotoluene | ND | 2.0 | 0.57 | ug/l | |
| 84-74-2Di-n-butyl phthalateND2.0 0.60 ug/l 117-84-0Di-n-octyl phthalateND 2.0 0.58 ug/l 84-66-2Diethyl phthalateND 2.0 0.40 ug/l 131-11-3Dimethyl phthalateND 2.0 0.33 ug/l 117-81-7bis(2-Ethylhexyl)phthalateND 2.0 0.67 ug/l 206-44-0FluorantheneND 1.0 0.25 ug/l 206-44-0FluorantheneND 1.0 0.25 ug/l 86-73-7FluoreneND 1.0 0.46 ug/l 118-74-1HexachlorobenzeneND 1.0 0.55 ug/l 87-68-3HexachlorocyclopentadieneND 1.0 0.42 ug/l 77-47-4HexachlorocyclopentadieneND 1.0 0.42 ug/l 193-39-5Indeno(1,2,3-cd)pyreneND 1.0 0.31 ug/l 91-20-3NaphthaleneND 1.0 0.33 ug/l 92-53NitrobenzeneND 2.0 0.43 ug/l 62-75-9n-NitrosodimethylamineND 2.0 0.48 ug/l 86-30-6N-NitrosodiphenylamineND 5.1 0.53 ug/l 85-01-8PhenanthreneND 1.0 0.34 ug/l 129-00-0PyreneND 1.0 0.34 ug/l 120-82-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.35 ug/l 12 | 91-94-1 | 3,3'-Dichlorobenzidine | ND | 2.0 | 1.2 | ug/l | |
| 117-84-0Di-n-octyl phthalateND2.0 0.58 ug/l 84-66-2Diethyl phthalateND2.0 0.40 ug/l 131-11-3Dimethyl phthalateND 2.0 0.33 ug/l 117-81-7bis(2-Ethylhexyl)phthalateND 2.0 0.67 ug/l 206-44-0FluorantheneND 1.0 0.25 ug/l 86-73-7FluoreneND 1.0 0.46 ug/l 118-74-1HexachlorobenzeneND 1.0 0.46 ug/l 87-68-3HexachlorobutadieneND 1.0 0.42 ug/l 67-72-1HexachlorocyclopentadieneND 1.0 0.31 ug/l 93-39-5Indeno(1,2,3-cd)pyreneND 1.0 0.31 ug/l 91-20-3NaphthaleneND 1.0 0.33 ug/l 98-95-3NitrobenzeneND 2.0 0.43 ug/l 62-75-9n-NitrosodimethylamineND 2.0 0.48 ug/l 86-30-6N-NitrosodiphenylamineND 2.0 0.48 ug/l 85-01-8PhenanthreneND 1.0 0.37 ug/l 129-09-0PyreneND 1.0 0.34 ug/l 129-09-1I,2,4-TrichlorobenzeneND 1.0 0.35 ug/l 129-09-2PhenathreneND 1.0 0.35 ug/l 129-09-3PyreneND 1.0 0.35 ug/l 120-09-4Pyrene </td <td>53-70-3</td> <td>Dibenzo(a,h)anthracene</td> <td>ND</td> <td>1.0</td> <td>0.56</td> <td>ug/l</td> <td></td> | 53-70-3 | Dibenzo(a,h)anthracene | ND | 1.0 | 0.56 | ug/l | |
| 84-66-2Diethyl phthalateND 2.0 0.40 ug/l 131-11-3Dimethyl phthalateND 2.0 0.33 ug/l 117-81-7bis(2-Ethylhexyl)phthalateND 2.0 0.67 ug/l 206-44-0FluorantheneND 1.0 0.25 ug/l 86-73-7FluoreneND 1.0 0.46 ug/l 118-74-1HexachlorobenzeneND 1.0 0.46 ug/l 87-68-3HexachlorobutadieneND 1.0 0.18 ug/l 77-47-4HexachlorocyclopentadieneND 1.0 0.42 ug/l 67-72-1HexachlorocthaneND 1.0 0.31 ug/l 93-39-5Indeno(1,2,3-cd)pyreneND 1.0 0.31 ug/l 98-95-3NitrobenzeneND 1.0 0.33 ug/l 98-95-3NitrobenzeneND 2.0 0.43 ug/l 62-75-9n-NitrosodimethylamineND 2.0 0.43 ug/l 62-64-7N-NitrosodiphenylamineND 2.0 0.48 ug/l 86-30-6N-NitrosodiphenylamineND 5.1 0.53 ug/l 129-00-0PyreneND 1.0 0.34 ug/l 129-82-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.34 ug/l 120-82-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.35 ug/l 120-82-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.35 ug/l | 84-74-2 | Di-n-butyl phthalate | ND | 2.0 | 0.60 | ug/l | |
| 131-11-3Dimethyl phthalateND 2.0 0.33 ug/l 117-81-7bis(2-Ethylhexyl)phthalateND 2.0 0.67 ug/l 206-44-0FluorantheneND 1.0 0.25 ug/l 86-73-7FluoreneND 1.0 0.46 ug/l 118-74-1HexachlorobenzeneND 1.0 0.46 ug/l 87-68-3HexachlorobutadieneND 1.0 0.18 ug/l 77-47-4HexachlorocyclopentadieneND 10 0.42 ug/l 67-72-1HexachloroethaneND 1.0 0.31 ug/l 93-39-5Indeno(1,2,3-cd)pyreneND 1.0 0.31 ug/l 91-20-3NaphthaleneND 1.0 0.33 ug/l 98-95-3NitrobenzeneND 2.0 0.43 ug/l 62-75-9n-NitrosodimethylamineND 2.0 0.443 ug/l 62-64-7N-NitrosodiphenylamineND 2.0 0.448 ug/l 86-30-6N-NitrosodiphenylamineND 5.1 0.53 ug/l 85-01-8PhenanthreneND 1.0 0.34 ug/l 129-00-0PyreneND 1.0 0.35 ug/l 129-00-0PyreneND 1.0 0.35 ug/l 120-82-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.35 ug/l 120-82-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.35 ug/l 120-8 | 117-84-0 | Di-n-octyl phthalate | ND | 2.0 | 0.58 | ug/l | |
| 117-81-7bis(2-Ethylhexyl)phthalateND 2.0 0.67 ug/l 206-44-0FluorantheneND 1.0 0.25 ug/l 86-73-7FluoreneND 1.0 0.46 ug/l 118-74-1HexachlorobenzeneND 1.0 0.46 ug/l 87-68-3HexachlorobutadieneND 1.0 0.18 ug/l 77-47-4HexachlorocyclopentadieneND 10 0.42 ug/l 67-72-1HexachloroethaneND 1.0 0.31 ug/l 93-39-5Indeno(1,2,3-cd)pyreneND 1.0 0.31 ug/l 91-20-3NaphthaleneND 1.0 0.33 ug/l 98-95-3NitrobenzeneND 2.0 0.43 ug/l 62-75-9n-NitrosodimethylamineND 2.0 0.44 ug/l 62-64-7N-Nitroso-di-n-propylamineND 2.0 0.44 ug/l 86-30-6N-NitrosodiphenylamineND 5.1 0.53 ug/l 85-01-8PhenanthreneND 1.0 0.37 ug/l 129-00-0PyreneND 1.0 0.35 ug/l 120-82-1 $1,2,4$ -TrichlorobenzeneND 1.0 0.35 ug/l 367-12-42-Fluorophenol 40% $14-110\%$ 4165-62-2Phenol-d5 29% $10-110\%$ | 84-66-2 | Diethyl phthalate | ND | 2.0 | 0.40 | ug/l | |
| 206-44-0FluorantheneND1.00.25ug/l86-73-7FluoreneND1.00.46ug/l118-74-1HexachlorobenzeneND1.00.55ug/l87-68-3HexachlorobutadieneND1.00.18ug/l77-47-4HexachlorocyclopentadieneND100.42ug/l67-72-1HexachloroethaneND5.10.29ug/l193-39-5Indeno(1,2,3-cd)pyreneND1.00.31ug/l91-20-3NaphthaleneND1.00.33ug/l92-53NitrobenzeneND2.00.60ug/l62-75-9n-NitrosodimethylamineND2.00.43ug/l62-64-7N-Nitrosodi-n-propylamineND2.00.48ug/l86-30-6N-NitrosodiphenylamineND5.10.53ug/l85-01-8PhenanthreneND1.00.34ug/l129-00-0PyreneND1.00.35ug/l120-82-11,2,4-TrichlorobenzeneND1.00.35ug/l367-12-42-Fluorophenol40%14-110%14-110%4165-62-2Phenol-d529%10-110%14-110% | 131-11-3 | Dimethyl phthalate | ND | 2.0 | 0.33 | ug/l | |
| 86-73-7 Fluorene ND 1.0 0.46 ug/l 118-74-1 Hexachlorobenzene ND 1.0 0.55 ug/l 87-68-3 Hexachlorobutadiene ND 1.0 0.18 ug/l 77-47-4 Hexachlorocyclopentadiene ND 10 0.42 ug/l 67-72-1 Hexachlorocyclopentadiene ND 5.1 0.29 ug/l 193-39-5 Indeno(1,2,3-cd)pyrene ND 1.0 0.31 ug/l 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.43 ug/l 62-64-7 N-Nitrosodiphenylamine ND 2.0 0.44 ug/l 86-30-6 N-Nitrosodiphenylamine ND 1.0 0.37 ug/l 85-01-8 Phenanthrene ND 1.0 0.34 ug/l 129-00-0 Pyrene ND 1.0 0.35 ug/l 120-82-1 1,2,4-Trich | 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 2.0 | 0.67 | ug/l | |
| 118-74-1 Hexachlorobenzene ND 1.0 0.55 ug/l 87-68-3 Hexachlorobutadiene ND 1.0 0.18 ug/l 77-47-4 Hexachlorocyclopentadiene ND 10 0.42 ug/l 67-72-1 Hexachloroethane ND 5.1 0.29 ug/l 193-39-5 Indeno(1,2,3-cd)pyrene ND 1.0 0.31 ug/l 78-59-1 Isophorone ND 2.0 0.60 ug/l 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.43 ug/l 62-64-7 N-Nitrosodiphenylamine ND 2.0 0.48 ug/l 86-30-6 N-Nitrosodiphenylamine ND 1.0 0.37 ug/l 120-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogat | 206-44-0 | Fluoranthene | ND | 1.0 | 0.25 | ug/l | |
| 87-68-3 Hexachlorobutadiene ND 1.0 0.18 ug/l 77-47-4 Hexachlorocyclopentadiene ND 10 0.42 ug/l 67-72-1 Hexachlorocyclopentadiene ND 5.1 0.29 ug/l 193-39-5 Indeno(1,2,3-cd)pyrene ND 1.0 0.31 ug/l 78-59-1 Isophorone ND 2.0 0.60 ug/l 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.44 ug/l 62-64-7 N-Nitroso-di-n-propylamine ND 2.0 0.48 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l 367-12-4 | 86-73-7 | Fluorene | ND | 1.0 | 0.46 | ug/l | |
| 77-47-4 Hexachlorocyclopentadiene ND 10 0.42 ug/l 67-72-1 Hexachloroethane ND 5.1 0.29 ug/l 193-39-5 Indeno(1,2,3-cd)pyrene ND 1.0 0.31 ug/l 78-59-1 Isophorone ND 2.0 0.60 ug/l 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.43 ug/l 621-64-7 N-Nitroso-di-n-propylamine ND 2.0 0.44 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 86-30-6 N-Nitrosodiphenylamine ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluoroph | 118-74-1 | Hexachlorobenzene | ND | 1.0 | 0.55 | ug/l | |
| 67-72-1 Hexachloroethane ND 5.1 0.29 ug/l 193-39-5 Indeno(1,2,3-cd)pyrene ND 1.0 0.31 ug/l 78-59-1 Isophorone ND 2.0 0.60 ug/l 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.443 ug/l 621-64-7 N-Nitroso-di-n-propylamine ND 2.0 0.448 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.35 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 87-68-3 | Hexachlorobutadiene | ND | 1.0 | 0.18 | ug/l | |
| 193-39-5 Indeno(1,2,3-cd)pyrene ND 1.0 0.31 ug/l 78-59-1 Isophorone ND 2.0 0.60 ug/l 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.47 ug/l 62-64-7 N-Nitroso-di-n-propylamine ND 2.0 0.48 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 77-47-4 | Hexachlorocyclopentadiene | ND | 10 | 0.42 | ug/l | |
| 78-59-1 Isophorone ND 2.0 0.60 ug/l 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.47 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.448 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.448 ug/l 62-64-7 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 86-30-6 N-Nitrosodiphenylamine ND 1.0 0.37 ug/l 85-01-8 Phenanthrene ND 1.0 0.34 ug/l 129-00-0 Pyrene ND 1.0 0.35 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 67-72-1 | Hexachloroethane | ND | 5.1 | 0.29 | ug/l | |
| 91-20-3 Naphthalene ND 1.0 0.33 ug/l 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.47 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.44 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.44 ug/l 62-64-7 N-Nitrosodiphenylamine ND 2.0 0.48 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 1.0 | 0.31 | ug/l | |
| 98-95-3 Nitrobenzene ND 2.0 0.43 ug/l 62-75-9 n-Nitrosodimethylamine ND 2.0 0.47 ug/l 621-64-7 N-Nitroso-di-n-propylamine ND 2.0 0.48 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 78-59-1 | Isophorone | ND | 2.0 | 0.60 | ug/l | |
| 62-75-9 n-Nitrosodimethylamine ND 2.0 0.47 ug/l 621-64-7 N-Nitroso-di-n-propylamine ND 2.0 0.48 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run# 1 Run# 2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 91-20-3 | Naphthalene | ND | 1.0 | 0.33 | ug/l | |
| 621-64-7 N-Nitroso-di-n-propylamine ND 2.0 0.48 ug/l 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 98-95-3 | Nitrobenzene | ND | 2.0 | 0.43 | ug/l | |
| 86-30-6 N-Nitrosodiphenylamine ND 5.1 0.53 ug/l 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 62-75-9 | n-Nitrosodimethylamine | ND | 2.0 | 0.47 | ug/l | |
| 85-01-8 Phenanthrene ND 1.0 0.37 ug/l 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 621-64-7 | N-Nitroso-di-n-propylamine | ND | 2.0 | 0.48 | ug/l | |
| 129-00-0 Pyrene ND 1.0 0.34 ug/l 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run#1 Run#2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 86-30-6 | N-Nitrosodiphenylamine | ND | 5.1 | 0.53 | ug/l | |
| 120-82-1 1,2,4-Trichlorobenzene ND 1.0 0.35 ug/l CAS No. Surrogate Recoveries Run# 1 Run# 2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 85-01-8 | Phenanthrene | ND | 1.0 | 0.37 | ug/l | |
| CAS No. Surrogate Recoveries Run# 1 Run# 2 Limits 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 129-00-0 | Pyrene | ND | 1.0 | 0.34 | ug/l | |
| 367-12-4 2-Fluorophenol 40% 14-110% 4165-62-2 Phenol-d5 29% 10-110% | 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.35 | ug/l | |
| 4165-62-2 Phenol-d5 29% 10-110% | CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| | 367-12-4 | | 40% | | 14-1 | 10% | |
| 119.70 C 2.4 C Tuil as we also and 790/ 20.1450/ | 4165-62-2 | Phenol-d5 | 29% | | 10-1 | 10% | |
| 118-79-6 2,4,6-Tribromophenol 78% 38-145% | 118-79-6 | 2,4,6-Tribromophenol | 78% | | 38-1 | 45% | |
| 4165-60-0 Nitrobenzene-d5 63% 33-136% | 4165-60-0 | Nitrobenzene-d5 | 63% | | 33-1 | 36% | |
| 321-60-8 2-Fluorobiphenyl 60% 35-127% | 321-60-8 | 2-Fluorobiphenyl | 60% | | 35-1 | 27% | |
| 1718-51-0 Terphenyl-d14 57% 11-139% | 1718-51-0 | | 57% | | 11-1 | 39% | |

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 2

SGS

5.0 ml

| Lab Sam Matrix: Method: Project: | AQ - G EPA 60 | 4061-3 - Ground Water A 608 EPA 608 2-2024 Cropsey Avenue, Brooklyn, NY | | | Date Sampled:02/10/16Date Received:02/10/16Percent Solids:n/a | | |
|---|-----------------------------|--|--------------------------|-----------------|---|-----------------------|----------------------------|
| | | | | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | File ID 6G31855.D | DF 1 | Analyzed 02/12/16 | By YD | Prep Date 02/12/16 | Prep Batch OP91166 | Analytical Batch G6G929 |
| Run #1 Run #2 | | DF 1 1 | • | v | 1 | 1 | • |

Report of Analysis

Pesticide PPL List

990 ml

Run #2

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|------------------------------|-----------------|--------|---------|-------|---|
| 309-00-2 | Aldrin | ND | 0.010 | 0.0017 | ug/l | |
| 319-84-6 | alpha-BHC | ND | 0.010 | 0.0013 | ug/l | |
| 319-85-7 | beta-BHC | ND | 0.010 | 0.0031 | ug/l | |
| 319-86-8 | delta-BHC | ND | 0.010 | 0.0016 | ug/l | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.010 | 0.00086 | ug/l | |
| 12789-03-6 | Chlordane | ND | 0.10 | 0.033 | ug/l | |
| 60-57-1 | Dieldrin | ND | 0.010 | 0.00086 | ug/l | |
| 72-54-8 | 4,4'-DDD | ND | 0.010 | 0.0012 | ug/l | |
| 72-55-9 | 4,4'-DDE | ND | 0.010 | 0.00084 | ug/l | |
| 50-29-3 | 4,4'-DDT | ND | 0.010 | 0.0025 | ug/l | |
| 72-20-8 | Endrin | ND | 0.010 | 0.0015 | ug/l | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.010 | 0.0023 | ug/l | |
| 7421-93-4 | Endrin aldehyde ^a | ND | 0.010 | 0.0032 | ug/l | |
| 959-98-8 | Endosulfan-I | ND | 0.010 | 0.0011 | ug/l | |
| 33213-65-9 | Endosulfan-II | ND | 0.010 | 0.0016 | ug/l | |
| 76-44-8 | Heptachlor | ND | 0.010 | 0.0013 | ug/l | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.010 | 0.00074 | ug/l | |
| 72-43-5 | Methoxychlor | ND | 0.010 | 0.0034 | ug/l | |
| 8001-35-2 | Toxaphene | ND | 0.13 | 0.047 | ug/l | |
| 12674-11-2 | Aroclor 1016 | ND ^b | 0.25 | 0.047 | ug/l | |
| 11104-28-2 | Aroclor 1221 | ND ^b | 0.25 | 0.24 | ug/l | |
| 11141-16-5 | Aroclor 1232 | ND ^b | 0.25 | 0.20 | ug/l | |
| 53469-21-9 | Aroclor 1242 | ND ^b | 0.25 | 0.083 | ug/l | |
| 12672-29-6 | Aroclor 1248 | ND ^b | 0.25 | 0.077 | ug/l | |
| 11097-69-1 | Aroclor 1254 | ND ^b | 0.25 | 0.055 | ug/l | |
| 11096-82-5 | Aroclor 1260 | ND ^b | 0.25 | 0.059 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 877-09-8 | Tetrachloro-m-xylene | 74% | 75% | 21-13 | 37% | |
| 877-09-8 | Tetrachloro-m-xylene | 75% | 84% | 21-13 | 87% | |
| 2051-24-3 | Decachlorobiphenyl | 46% | 39% | 10-12 | 21% | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



19 of 963

JC14061

4.3

Page 1 of 2

Report of Analysis

Client Sample ID:MW-2Lab Sample ID:JC14061-3Date Sampled:02/10/16Matrix:AQ - Ground WaterDate Received:02/10/16Method:EPA 608EPA 608Percent Solids:n/aProject:2002-2024 Cropsey Avenue, Brooklyn, NYNY

Pesticide PPL List

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 2051-24-3 | Decachlorobiphenyl | 54% | 40% | 10-121% |

(a) This compound outside control limits biased high in the associated BS.

(b) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

SGS

20 of 963

ACCUTEST JC14061

4.3 4

| Report of Analysis | | | | | | | | | |
|---|---------------------------------|--------------------------|--------------------------|-----------|---------------------------|-----------------------|-----------------------------|--|--|
| Client Sample ID:MW-2Lab Sample ID:JC14061-3Matrix:AQ - Ground WaterMethod:SW846 8151SW846 8151SW846 8151/3510CProject:2002-2024 Cropsey Avenue, Brooklyn, NY | | | | | | | | | |
| Run #1 Run #2 | File ID OA114320.D | DF 1 | Analyzed 02/16/16 | By VDT | Prep Date 02/12/16 | Prep Batch OP91147 | Analytical Batch GOA3955 | | |
| Run #1 Run #2 | Initial Volume 980 ml | Final \ 10.0 m | Volume hl | | | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|----------------------|------------------------|----------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 0.51 0.10 0.10 | 0.29 0.056 0.057 | ug/l ug/l ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2,4-DCAA | 107% 80% | | 39-1 39-1 | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



21 of 963 ACCUTEST JC14061

| Client Sample ID: | MW-2 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC14061-3 | Date Sampled: | 02/10/16 |
| Matrix: | AQ - Ground Water | Date Received: | 02/10/16 |
| | | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Total Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|--------|------|-------|----|----------|-------------|--------------------------|--------------------------|
| Antimony | < 6.0 | 6.0 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Arsenic | < 3.0 | 3.0 | ug/l | 1 | 02/12/16 | 02/12/16 BS | SW846 6010C ² | SW846 3010A ³ |
| Beryllium | < 1.0 | 1.0 | ug/l | 1 | 02/12/16 | 02/12/16 BS | SW846 6010C ² | SW846 3010A ³ |
| Cadmium | < 3.0 | 3.0 | ug/l | 1 | 02/12/16 | 02/12/16 BS | SW846 6010C ² | SW846 3010A ³ |
| Chromium | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 BS | SW846 6010C ² | SW846 3010A ³ |
| Copper | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 BS | SW846 6010C ² | SW846 3010A ³ |
| Lead | < 3.0 | 3.0 | ug/l | 1 | 02/12/16 | 02/12/16 вѕ | SW846 6010C ² | SW846 3010A ³ |
| Mercury | < 0.20 | 0.20 | ug/l | 1 | 02/12/16 | 02/12/16 VM | SW846 7470A ¹ | SW846 7470A ⁴ |
| Nickel | 18.5 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вѕ | SW846 6010C ² | SW846 3010A ³ |
| Selenium | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вѕ | SW846 6010C ² | SW846 3010A ³ |
| Silver | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вѕ | SW846 6010C ² | SW846 3010A ³ |
| Thallium | < 2.0 | 2.0 | ug/l | 1 | 02/12/16 | 02/12/16 вѕ | SW846 6010C ² | SW846 3010A ³ |
| Zinc | < 20 | 20 | ug/l | 1 | 02/12/16 | 02/12/16 BS | SW846 6010C ² | SW846 3010A ³ |

(1) Instrument QC Batch: MA38721

(2) Instrument QC Batch: MA38722

(3) Prep QC Batch: MP91988

(4) Prep QC Batch: MP91993

Page 1 of 1





| Report of Analysis | | | | | | | | | | |
|---|---------------------------|--|----------------------------|-----------|---------------------------------------|--|-----------------------------------|--|--|--|
| Client Sar Lab Sam Matrix: Method: Project: | ple ID: | MW-3 IC14061-4 AQ - Ground V EPA 624 2002-2024 Cro | Water opsey Avenue, Bro | ooklyn, N | · · · · · · · · · · · · · · · · · · · | Date Sampled: Date Received: Percent Solids: | | | | |
| Run #1 ^a Run #2 | File ID U203281 | DF .D 1 | Analyzed 02/11/16 | By NH | Prep Date n/a | Prep Batch n/a | Analytical Batch VU9349 | | | |
| Run #1 | Purge V 5.0 ml | olume | | | | | | | | |

Run #2

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | 0.51 | 1.0 | 0.091 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 0.24 | 1.0 | 0.12 | ug/l | J |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 11.9 | 1.0 | 0.14 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



SGS

E = Indicates value exceeds calibration range

| Client Samp Lab Sample Matrix: Method: Project: | | MW-3 JC14061-4 AQ - Ground Water EPA 624 2002-2024 Cropsey Ave | enue, Brook | lyn, NY | | Date | Sampled: Received: nt Solids: | 02/10/16 02/10/16 n/a |
|---|----------------------|--|-------------|---------|---------|-------|-------------------------------------|-----------------------------|
| VOA PPL List | | | | | | | | |
| CAS No. | Compound | | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichl | Trichloroethene | | 1.0 | 0.12 | ug/l | J | |
| 75-69-4 | Trichl | Trichlorofluoromethane | | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl chloride | | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylenes (total) | | ND | 1.0 | 0.22 | ug/l | | |
| CAS No. | Surrogate Recoveries | | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Di | chloroethane-D4 (SUR) | 110% | 72-125% | | | | |
| 2037-26-5 | Toluer | ene-D8 (SUR) 101% | | | 78-119% | | | |
| 460-00-4 | 4-Broi | nofluorobenzene (SUR) | 102% | 74-115% | | | | |
| 1868-53-7 | Dibro | mofluoromethane (S) | 104% | 79-120% | | | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4.4

Page 2 of 2



| Report of Analysis Page 1 | | | | | | | | |
|--|---------------------------|--|----------------------------|-----------------|-------------------------|---|-----------------------------------|--|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: J A H | MW-4S C14061-5 AQ - Ground Y EPA 624 2002-2024 Cro | Water opsey Avenue, Bro | ooklyn, N |] | Date Sampled: 0. Date Received: 0. Percent Solids: n. | | |
| Run #1 ^a Run #2 | File ID U203282 | DF .D 1 | Analyzed 02/11/16 | By NH | Prep Date n/a | Prep Batch n/a | Analytical Batch VU9349 | |
| Run #1 | Purge Vo 5.0 ml | olume | | | | | | |

Run #2

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | 0.56 | 1.0 | 0.091 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 0.29 | 1.0 | 0.14 | ug/l | J |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $J = \ Indicates \ an \ estimated \ value$

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



SGS

E = Indicates value exceeds calibration range

| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | | MW-4S JC14061-5 AQ - Ground Water EPA 624 2002-2024 Cropsey Av | enue, Brook | Date | Sampled: Received: ent Solids: | 02/10/16 02/10/16 n/a | | |
|---|---------|--|-------------|--------|--------------------------------------|-----------------------------|---|--|
| VOA PPL I | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichlo | proethene | ND | 1.0 | 0.12 | ug/l | | |
| 75-69-4 | Trichle | orofluoromethane | ND | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl o | chloride | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylene | es (total) | ND | 1.0 | 0.22 | ug/l | | |
| CAS No. | Surrog | gate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Di | chloroethane-D4 (SUR) | 110% | | 72-1 | 25% | | |
| 2037-26-5 | | | 101% | | 78-1 | 19% | | |
| 460-00-4 | 4-Bron | nofluorobenzene (SUR) | 103% | | 74-1 | 15% | | |
| 1868-53-7 | Dibror | nofluoromethane (S) | 105% | | 79-1 | 20% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



4.5

Page 2 of 2



| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|--|---------------------------|--|---------------------------|-----------------|-------------------------|--|-----------------------------------|
| Client Sar Lab Samp Matrix: Method: Project: | ble ID: J | MW-4I C14061-6 AQ - Ground W EPA 624 2002-2024 Cro | Vater psey Avenue, Bro | ooklyn, N |] | Date Sampled: 0 Date Received: 0 Percent Solids: n | |
| Run #1 ^a Run #2 | File ID U203283 | DF .D 1 | Analyzed 02/11/16 | By NH | Prep Date n/a | Prep Batch n/a | Analytical Batch VU9349 |
| Run #1 | Purge Vo 5.0 ml | olume | | | | | |

Run #2

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | 0.52 | 1.0 | 0.091 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 0.26 | 1.0 | 0.14 | ug/l | J |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



SGS

E = Indicates value exceeds calibration range

| Client Samp Lab Sample Matrix: Method: Project: | | venue, Brool | Date | Sampled: Received: ent Solids: | 02/10/16 02/10/16 n/a | | |
|---|-----------------------------|--------------|--------|--------------------------------------|-----------------------------|---|--|
| VOA PPL I | list | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.12 | ug/l | | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylenes (total) | ND | 1.0 | 0.22 | ug/l | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 (SUR) | 108% | | 72-1 | 25% | | |
| 2037-26-5 | Toluene-D8 (SUR) | 101% | | 78-1 | 19% | | |
| 460-00-4 | 4-Bromofluorobenzene (SUR) | 102% | | 74-1 | 15% | | |
| 1868-53-7 | Dibromofluoromethane (S) | 104% | | 79-1 | 20% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2



| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|--|------------------------------|----------------|----------------------------|-----------------|-------------------------|--|------------------------------|
| Client Sar Lab Samp Matrix: Method: Project: | Ile ID: JC14 AQ - EPA | | c Water sey Avenue, Bro | ooklyn, N | Y | Date Sampled: Date Received: Percent Solids: | |
| Run #1 ^a Run #2 | File ID U203277.D | DF 1 | Analyzed 02/11/16 | By NH | Prep Date n/a | Prep Batcl n/a | n Analytical Batch VU9349 |
| Run #1 Run #2 | Purge Volun 5.0 ml | ne | | | | | |

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.091 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.14 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



29 of 963

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

1868-53-7

| Client Samp Lab Sample Matrix: Method: Project: | | FIELD/FB JC14061-7 AQ - Field Blank Water EPA 624 2002-2024 Cropsey Ave | | lyn, NY | | Date | Sampled: Received: ent Solids: | 02/10/16 02/10/16 n/a |
|---|-----------------|---|----------------------|-------------------------------|------------------------------|------------------------------|--------------------------------------|-----------------------------|
| VOA PPL I | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | |
| 79-01-6 75-69-4 75-01-4 1330-20-7 | Trichl Vinyl | Trichloroethene Trichlorofluoromethane Vinyl chloride Xylenes (total) | | 1.0 2.0 1.0 1.0 | 0.12 0.20 0.13 0.22 | ug/l ug/l ug/l ug/l | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 2037-26-5 460-00-4 | Toluer | ichloroethane-D4 (SUR) ne-D8 (SUR) nofluorobenzene (SUR) | 109% 102% 103% | 72-125% 78-119% 74-115% | | 19% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

103%

79-120%

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



4.7



| Client San Lab Samj Matrix: Method: Project: | ple ID: JC14 AQ EPA | LD/FB 1061-7 - Field Blank 625 EPA 2-2024 Crops | | ooklyn, NY | Da Pe | ate Sampled: (ate Received: (prcent Solids: 1 | |
|--|--------------------------------|---|--------------------------|------------|---------------------------|---|-----------------------------------|
| Run #1 Run #2 | File ID F154588.D | DF 1 | Analyzed 02/12/16 | By SD | Prep Date 02/11/16 | Prep Batch OP91141 | Analytical Batch EF6502 |
| Run #1 | Initial Volur 975 ml | ne Final V | olume | | | | |

Report of Analysis

Run #2

1.0 ml

ABN PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|-----|------|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 5.1 | 0.98 | ug/l | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 5.1 | 1.2 | ug/l | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 5.1 | 1.6 | ug/l | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 5.1 | 1.7 | ug/l | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 5.1 | 0.91 | ug/l | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 5.1 | 0.74 | ug/l | |
| 88-75-5 | 2-Nitrophenol | ND | 2.1 | 1.9 | ug/l | |
| 100-02-7 | 4-Nitrophenol | ND | 10 | 0.86 | ug/l | |
| 87-86-5 | Pentachlorophenol | ND | 5.1 | 1.9 | ug/l | |
| 108-95-2 | Phenol | ND | 2.1 | 0.51 | ug/l | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 2.1 | 1.3 | ug/l | |
| 83-32-9 | Acenaphthene | ND | 1.0 | 0.36 | ug/l | |
| 208-96-8 | Acenaphthylene | ND | 1.0 | 0.39 | ug/l | |
| 120-12-7 | Anthracene | ND | 1.0 | 0.41 | ug/l | |
| 92-87-5 | Benzidine | ND | 21 | 0.29 | ug/l | |
| 56-55-3 | Benzo(a)anthracene | ND | 1.0 | 0.37 | ug/l | |
| 50-32-8 | Benzo(a)pyrene | ND | 1.0 | 0.38 | ug/l | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 1.0 | 0.61 | ug/l | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 1.0 | 0.43 | ug/l | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 1.0 | 0.43 | ug/l | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 2.1 | 0.31 | ug/l | |
| 85-68-7 | Butyl benzyl phthalate | ND | 2.1 | 0.61 | ug/l | |
| 91-58-7 | 2-Chloronaphthalene | ND | 2.1 | 1.0 | ug/l | |
| 106-47-8 | 4-Chloroaniline | ND | 2.1 | 0.41 | ug/l | |
| 218-01-9 | Chrysene | ND | 1.0 | 0.26 | ug/l | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 2.1 | 0.67 | ug/l | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 2.1 | 0.54 | ug/l | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 2.1 | 0.76 | ug/l | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 2.1 | 0.44 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 2.1 | 0.21 | ug/l | |
| 122-66-7 | 1,2-Diphenylhydrazine | ND | 2.1 | 0.46 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 2.1 | 0.16 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2

JC14061

SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

Client Sample ID:FIELD/FBLab Sample ID:JC14061-7Date Sampled:02/10/16Matrix:AQ - Field Blank WaterDate Received:02/10/16Method:EPA 625EPA 625Percent Solids:n/aProject:2002-2024 Cropsey Avenue, Brooklyn, NYNYNY

ABN PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|----------------------------|--------|--------|------|-------|---|
| 106-46-7 | 1,4-Dichlorobenzene | ND | 2.1 | 0.18 | ug/l | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 2.1 | 0.88 | ug/l | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 2.1 | 0.57 | ug/l | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 2.1 | 1.3 | ug/l | |
| 53-70-3 | Dibenzo(a, h)anthracene | ND | 1.0 | 0.56 | ug/l | |
| 84-74-2 | Di-n-butyl phthalate | ND | 2.1 | 0.61 | ug/l | |
| 117-84-0 | Di-n-octyl phthalate | ND | 2.1 | 0.58 | ug/l | |
| 84-66-2 | Diethyl phthalate | ND | 2.1 | 0.40 | ug/l | |
| 131-11-3 | Dimethyl phthalate | ND | 2.1 | 0.34 | ug/l | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 2.1 | 0.68 | ug/l | |
| 206-44-0 | Fluoranthene | ND | 1.0 | 0.26 | ug/l | |
| 86-73-7 | Fluorene | ND | 1.0 | 0.46 | ug/l | |
| 118-74-1 | Hexachlorobenzene | ND | 1.0 | 0.55 | ug/l | |
| 87-68-3 | Hexachlorobutadiene | ND | 1.0 | 0.18 | ug/l | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 10 | 0.42 | ug/l | |
| 67-72-1 | Hexachloroethane | ND | 5.1 | 0.29 | ug/l | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 1.0 | 0.31 | ug/l | |
| 78-59-1 | Isophorone | ND | 2.1 | 0.60 | ug/l | |
| 91-20-3 | Naphthalene | ND | 1.0 | 0.33 | ug/l | |
| 98-95-3 | Nitrobenzene | ND | 2.1 | 0.43 | ug/l | |
| 62-75-9 | n-Nitrosodimethylamine | ND | 2.1 | 0.47 | ug/l | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 2.1 | 0.48 | ug/l | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 5.1 | 0.53 | ug/l | |
| 85-01-8 | Phenanthrene | ND | 1.0 | 0.37 | ug/l | |
| 129-00-0 | Pyrene | ND | 1.0 | 0.34 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.35 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 367-12-4 | 2-Fluorophenol | 41% | | 14-1 | 10% | |
| 4165-62-2 | Phenol-d5 | 29% | | 10-1 | 10% | |
| 118-79-6 | 2,4,6-Tribromophenol | 76% | | 38-1 | 45% | |
| 4165-60-0 | Nitrobenzene-d5 | 59% | | 33-1 | 36% | |
| 321-60-8 | 2-Fluorobiphenyl | 58% | | 35-1 | 27% | |
| 1718-51-0 | Terphenyl-d14 | 70% | | 11-1 | 39% | |

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 2 of 2



5.0 ml

| Client Sar Lab Samj Matrix: Method: Project: | ple ID: JC1406 AQ - F EPA 60 | il-7 ield Blan)8 EPA | | ooklyn N | Da Pe | 1 | 2/10/16 2/10/16 /a |
|--|------------------------------------|-----------------------------|----------------------|----------|----------------------|--------------------|--------------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 Run #2 | 6G31856.D XX185092.D | 1 1 | 02/12/16 02/12/16 | YD KM | 02/12/16 02/12/16 | OP91166 OP91165 | G6G929 GXX5601 |
| Run #1 | Initial Volume 1000 ml | Final 5.0 ml | Volume | | | | |

Report of Analysis

Pesticide PPL List

970 ml

Run #2

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|------------------------------|-----------------|--------|---------|-------|---|
| 309-00-2 | Aldrin | ND | 0.010 | 0.0017 | ug/l | |
| 319-84-6 | alpha-BHC | ND | 0.010 | 0.0013 | ug/l | |
| 319-85-7 | beta-BHC | ND | 0.010 | 0.0031 | ug/l | |
| 319-86-8 | delta-BHC | ND | 0.010 | 0.0016 | ug/l | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.010 | 0.00086 | ug/l | |
| 12789-03-6 | Chlordane | ND | 0.10 | 0.033 | ug/l | |
| 60-57-1 | Dieldrin | ND | 0.010 | 0.00086 | ug/l | |
| 72-54-8 | 4,4'-DDD | ND | 0.010 | 0.0012 | ug/l | |
| 72-55-9 | 4,4'-DDE | ND | 0.010 | 0.00084 | ug/l | |
| 50-29-3 | 4,4'-DDT | ND | 0.010 | 0.0025 | ug/l | |
| 72-20-8 | Endrin | ND | 0.010 | 0.0015 | ug/l | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.010 | 0.0023 | ug/l | |
| 7421-93-4 | Endrin aldehyde ^a | ND | 0.010 | 0.0032 | ug/l | |
| 959-98-8 | Endosulfan-I | ND | 0.010 | 0.0011 | ug/l | |
| 33213-65-9 | Endosulfan-II | ND | 0.010 | 0.0016 | ug/l | |
| 76-44-8 | Heptachlor | ND | 0.010 | 0.0013 | ug/l | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.010 | 0.00074 | ug/l | |
| 72-43-5 | Methoxychlor | ND | 0.010 | 0.0034 | ug/l | |
| 8001-35-2 | Toxaphene | ND | 0.13 | 0.047 | ug/l | |
| 12674-11-2 | Aroclor 1016 | ND ^b | 0.26 | 0.048 | ug/l | |
| 11104-28-2 | Aroclor 1221 | ND ^b | 0.26 | 0.24 | ug/l | |
| 11141-16-5 | Aroclor 1232 | ND ^b | 0.26 | 0.20 | ug/l | |
| 53469-21-9 | Aroclor 1242 | ND ^b | 0.26 | 0.085 | ug/l | |
| 12672-29-6 | Aroclor 1248 | ND ^b | 0.26 | 0.079 | ug/l | |
| 11097-69-1 | Aroclor 1254 | ND ^b | 0.26 | 0.056 | ug/l | |
| 11096-82-5 | Aroclor 1260 | ND ^b | 0.26 | 0.060 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 877-09-8 | Tetrachloro-m-xylene | 107% | 113% | 21-13 | 37% | |
| 877-09-8 | Tetrachloro-m-xylene | 108% | 119% | 21-13 | 87% | |
| 2051-24-3 | Decachlorobiphenyl | 83% | 70% | 10-12 | 21% | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 1 of 2



Report of Analysis

Client Sample ID:FIELD/FBLab Sample ID:JC14061-7Date Sampled:02/10/16Matrix:AQ - Field Blank WaterDate Received:02/10/16Method:EPA 608EPA 608Percent Solids:n/aProject:2002-2024 Cropsey Avenue, Brooklyn, NYYClient Solids:N/a

Pesticide PPL List

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 2051-24-3 | Decachlorobiphenyl | 97% | 72% | 10-121% |

(a) This compound outside control limits biased high in the associated BS.

(b) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

SGS

34 of 963 ACCUTEST

Page 2 of 2

4.7



| | Report of Analysis | | | | | | | | | |
|----------------------|--------------------|------------|-----------------|-----------|-----------|-------------------|------------------|--|--|--|
| Client Sa Lab Sam | - | | | | Da | ate Sampled: 02 | 2/10/16 | | | |
| Matrix: | - | Field Blan | k Water | | | ate Received: 02 | 2/10/16 | | | |
| Method: | SW84 | 6 8151 S | W846 8151/351 | 0C | Pe | ercent Solids: n/ | a | | | |
| Project: | 2002- | 2024 Crop | sey Avenue, Bro | oklyn, NY | ζ. | | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | | | |
| Run #1 | OA114321.D | 1 | 02/16/16 | VDT | 02/12/16 | OP91147 | GOA3955 | | | |
| Run #2 | | | | | | | | | | |
| | Initial Volum | e Final V | Volume | | | | | | | |
| Run #1 | 960 ml | 10.0 m | 1 | | | | | | | |
| Run #2 | | | | | | | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|----------------------|------------------------|----------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 0.52 0.10 0.10 | 0.30 0.057 0.058 | ug/l ug/l ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 19719-28-9 | 2,4-DCAA | 107% | | 39-1 | 59% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



35 of 963 ACCUTEST JC14061

4.7

4

 Client Sample ID:
 FIELD/FB

 Lab Sample ID:
 JC14061-7

 Matrix:
 AQ - Field Blank Water

 Project:
 2002-2024 Cropsey Avenue, Brooklyn, NY

Total Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|--------|------|-------|----|----------|-------------|--------------------------|--------------------------|
| Antimony | < 6.0 | 6.0 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Arsenic | < 3.0 | 3.0 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Beryllium | < 1.0 | 1.0 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Cadmium | < 3.0 | 3.0 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Chromium | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C 2 | SW846 3010A ³ |
| Copper | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Lead | < 3.0 | 3.0 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Mercury | < 0.20 | 0.20 | ug/l | 1 | 02/12/16 | 02/12/16 VM | SW846 7470A ¹ | SW846 7470A ⁴ |
| Nickel | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Selenium | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C ² | SW846 3010A ³ |
| Silver | < 10 | 10 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C 2 | SW846 3010A ³ |
| Thallium | < 2.0 | 2.0 | ug/l | 1 | 02/12/16 | 02/12/16 вз | SW846 6010C 2 | SW846 3010A ³ |
| Zinc | < 20 | 20 | ug/l | 1 | 02/12/16 | 02/12/16 вѕ | SW846 6010C 2 | SW846 3010A ³ |

(1) Instrument QC Batch: MA38721

(2) Instrument QC Batch: MA38722

(3) Prep QC Batch: MP91988

(4) Prep QC Batch: MP91993







| Report of AnalysisPage 1Client Sample ID:MW-ALab Sample ID:JC14061-8Matrix:AQ - Ground WaterMethod:EPA 624Project:2002-2024 Cropsey Avenue, Brooklyn, NY | | | | | | | | | | |
|--|---------------------|------|--|--|--|--|--|--|--|--|
| | | | | | | | | | | |
| Run #1 | Purge Vol 5.0 ml | lume | | | | | | | | |

Run #2

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.091 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.14 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



37 of 963

E = Indicates value exceeds calibration range

| Client Samp Lab Sample Matrix: Method: Project: | ID: JC14061-8 AQ - Ground W EPA 624 | Vater psey Avenue, Broo | klyn, NY | | Date | Sampled: Received: ent Solids: | 02/10/16 02/10/16 n/a |
|---|---|----------------------------|----------|------|-------|--------------------------------------|-----------------------------|
| VOA PPL I | List | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.12 | ug/l | | |
| 75-69-4 | Trichlorofluoromethan | e ND | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylenes (total) | ND | 1.0 | 0.22 | ug/l | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 4 (SUR) 111% | | 72-1 | 25% | | |
| 2037-26-5 | Toluene-D8 (SUR) | 101% | | 78-1 | 19% | | |
| 460-00-4 | 4-Bromofluorobenzene | e (SUR) 105% | | 74-1 | 15% | | |
| 1868-53-7 | Dibromofluoromethan | e (S) 106% | | 79-1 | 20% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



Page 2 of 2





Section 4

Sample Results

Report of Analysis





| Lab Sample ID:JC16575-1Matrix:AQ - Ground WaterMethod:EPA 624Project:2002-2024 Cropsey Avenue, Brooklyn, NY | | | | | Date Sampled:03/18/16Date Received:03/18/16Percent Solids:n/a | | |
|---|-----------|-----|----------|----|---|------------|------------------|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 ^a | N253486.D | 10 | 03/19/16 | wo | n/a | n/a | VN10722 |
| Run #2 ^a | N253487.D | 100 | 03/19/16 | WO | n/a | n/a | VN10722 |
| | | | | | | | |

Report of Analysis

Run #1 5.0 ml

Run #2 5.0 ml

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|-------------------|-----|------|-------|---|
| 107-02-8 | Acrolein | ND | 100 | 16 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 100 | 26 | ug/l | |
| 71-43-2 | Benzene | ND | 10 | 1.0 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 10 | 1.0 | ug/l | |
| 75-25-2 | Bromoform | ND | 10 | 1.7 | ug/l | |
| 74-83-9 | Bromomethane | ND | 10 | 5.7 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 10 | 0.96 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 10 | 0.93 | ug/l | |
| 75-00-3 | Chloroethane | ND | 10 | 2.1 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 50 | 5.0 | ug/l | |
| 67-66-3 | Chloroform | ND | 10 | 0.91 | ug/l | |
| 74-87-3 | Chloromethane | ND | 10 | 1.1 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 10 | 1.5 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 10 | 1.9 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 10 | 1.9 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 10 | 1.1 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 20 | 2.9 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 10 | 1.2 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 10 | 0.90 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 10 | 1.6 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 687 | 10 | 1.2 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 10 | 1.4 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 10 | 1.1 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 10 | 1.2 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 10 | 1.5 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 10 | 2.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 10 | 2.2 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 10 | 1.2 | ug/l | |
| 127-18-4 | Tetrachloroethene | 3490 ^b | 100 | 14 | ug/l | |
| 108-88-3 | Toluene | ND | 10 | 2.5 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 10 | 0.86 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 10 | 1.3 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 1 of 2

SGS

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

| Client Samp Lab Sample Matrix: Method: Project: | | venue, Brook | lyn, NY | | Date | Sampled: Received: ent Solids: | 03/18/16 03/18/16 n/a |
|---|--|-----------------------|----------------------|--------------------------|------------------------------|--------------------------------------|-----------------------------|
| VOA PPL I | ist | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 75-69-4 75-01-4 1330-20-7 | Trichloroethene Trichlorofluoromethane Vinyl chloride Xylenes (total) | 106 ND ND ND | 10 20 10 10 | 1.2 2.0 1.3 2.2 | ug/l ug/l ug/l ug/l | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | | |
| 17060-07-0 2037-26-5 460-00-4 | 1,2-Dichloroethane-D4 (SUR) Toluene-D8 (SUR) 4-Bromofluorobenzene (SUR) | 101% | 107% 100% 103% | 72-1 78-1 74-1 | 19% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

104%

102%

79-120%

(b) Result is from Run# 2

1868-53-7

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



4

| | | | r - | | | | |
|-----------|----------------|-----------|-----------------|----------|-----------|------------------|------------------|
| Client Sa | mple ID: MW-5 | | | | | | |
| Lab Sam | ple ID: JC1657 | 75-1 | | | Da | te Sampled: 03 | 8/18/16 |
| Matrix: | - AQ - G | bround Wa | ater | | Da | te Received: 03 | 8/18/16 |
| Method: | EPA 62 | 25 EPA | 625 | | Pe | rcent Solids: n/ | a |
| Project: | 2002-2 | 024 Crops | sey Avenue, Bro | oklyn, N | Y | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | F155736.D | 1 | 03/22/16 | SD | 03/21/16 | OP92311 | EF6551 |
| Run #2 | | | | | | | |
| | Initial Volume | Final V | olume | | | | |
| Run #1 | 970 ml | 1.0 ml | | | | | |

Report of Analysis

Run #2

ABN PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|-----|------|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 5.2 | 0.98 | ug/l | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 5.2 | 1.2 | ug/l | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 5.2 | 1.6 | ug/l | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 5.2 | 1.7 | ug/l | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 5.2 | 0.92 | ug/l | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 5.2 | 0.74 | ug/l | |
| 88-75-5 | 2-Nitrophenol | ND | 2.1 | 1.9 | ug/l | |
| 100-02-7 | 4-Nitrophenol | ND | 10 | 0.87 | ug/l | |
| 87-86-5 | Pentachlorophenol | ND | 5.2 | 2.0 | ug/l | |
| 108-95-2 | Phenol | ND | 2.1 | 0.51 | ug/l | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 2.1 | 1.3 | ug/l | |
| 83-32-9 | Acenaphthene | ND | 1.0 | 0.36 | ug/l | |
| 208-96-8 | Acenaphthylene | ND | 1.0 | 0.39 | ug/l | |
| 120-12-7 | Anthracene | ND | 1.0 | 0.41 | ug/l | |
| 92-87-5 | Benzidine | ND | 21 | 0.29 | ug/l | |
| 56-55-3 | Benzo(a)anthracene | ND | 1.0 | 0.37 | ug/l | |
| 50-32-8 | Benzo(a)pyrene | ND | 1.0 | 0.38 | ug/l | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 1.0 | 0.61 | ug/l | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 1.0 | 0.44 | ug/l | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 1.0 | 0.43 | ug/l | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 2.1 | 0.31 | ug/l | |
| 85-68-7 | Butyl benzyl phthalate | ND | 2.1 | 0.61 | ug/l | |
| 91-58-7 | 2-Chloronaphthalene | ND | 2.1 | 1.0 | ug/l | |
| 106-47-8 | 4-Chloroaniline | ND | 2.1 | 0.41 | ug/l | |
| 218-01-9 | Chrysene | ND | 1.0 | 0.26 | ug/l | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 2.1 | 0.67 | ug/l | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 2.1 | 0.55 | ug/l | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 2.1 | 0.76 | ug/l | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 2.1 | 0.44 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 2.1 | 0.21 | ug/l | |
| 122-66-7 | 1,2-Diphenylhydrazine | ND | 2.1 | 0.47 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 2.1 | 0.16 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Page 1 of 2

11 of 1072

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| | N-5 | | |
|--------------------|--------------------------------------|-----------------|----------|
| Lab Sample ID: JC1 | 16575-1 | Date Sampled: | 03/18/16 |
| Matrix: AQ | - Ground Water | Date Received: | 03/18/16 |
| Method: EPA | A 625 EPA 625 | Percent Solids: | n/a |
| Project: 200 | 02-2024 Cropsey Avenue, Brooklyn, NY | | |

ABN PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|----------------------------|--------|--------|------|-------|---|
| 106-46-7 | 1,4-Dichlorobenzene | ND | 2.1 | 0.19 | ug/l | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 2.1 | 0.89 | ug/l | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 2.1 | 0.58 | ug/l | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 2.1 | 1.3 | ug/l | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 1.0 | 0.56 | ug/l | |
| 84-74-2 | Di-n-butyl phthalate | ND | 2.1 | 0.61 | ug/l | |
| 117-84-0 | Di-n-octyl phthalate | ND | 2.1 | 0.59 | ug/l | |
| 84-66-2 | Diethyl phthalate | ND | 2.1 | 0.40 | ug/l | |
| 131-11-3 | Dimethyl phthalate | ND | 2.1 | 0.34 | ug/l | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 2.1 | 0.68 | ug/l | |
| 206-44-0 | Fluoranthene | ND | 1.0 | 0.26 | ug/l | |
| 86-73-7 | Fluorene | ND | 1.0 | 0.46 | ug/l | |
| 118-74-1 | Hexachlorobenzene | ND | 1.0 | 0.55 | ug/l | |
| 87-68-3 | Hexachlorobutadiene | ND | 1.0 | 0.18 | ug/l | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 10 | 0.42 | ug/l | |
| 67-72-1 | Hexachloroethane | ND | 5.2 | 0.29 | ug/l | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 1.0 | 0.31 | ug/l | |
| 78-59-1 | Isophorone | ND | 2.1 | 0.61 | ug/l | |
| 91-20-3 | Naphthalene | ND | 1.0 | 0.33 | ug/l | |
| 98-95-3 | Nitrobenzene | ND | 2.1 | 0.43 | ug/l | |
| 62-75-9 | n-Nitrosodimethylamine | ND | 2.1 | 0.47 | ug/l | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 2.1 | 0.48 | ug/l | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 5.2 | 0.53 | ug/l | |
| 85-01-8 | Phenanthrene | ND | 1.0 | 0.37 | ug/l | |
| 129-00-0 | Pyrene | ND | 1.0 | 0.35 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.35 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 367-12-4 | 2-Fluorophenol | 40% | | 14-1 | 10% | |
| 4165-62-2 | Phenol-d5 | 28% | | 10-1 | 10% | |
| 118-79-6 | 2,4,6-Tribromophenol | 71% | | 38-1 | 45% | |
| 4165-60-0 | Nitrobenzene-d5 | 71% | | 33-1 | 36% | |
| 321-60-8 | 2-Fluorobiphenyl | 65% | | 35-1 | 27% | |
| 1718-51-0 | Terphenyl-d14 | 68% | | 11-1 | 39% | |
| | | | | | | |

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



4.1 **4**

Page 2 of 2



12 of 1072

5.0 ml

| Client Sa Lab Samj Matrix: Method: | AQ - G | 5-1 round Wa | | | Da | te Sampled: 03 te Received: 03 rcent Solids: n/ | |
|---|---|-----------------|----------|----|-----------|---|------------------|
| Project: | Project: 2002-2024 Cropsey Avenue, Brooklyn, NY | | | | Ŷ | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | 6G33312.D | 1 | 03/24/16 | BP | 03/22/16 | OP92352 | G6G969 |
| Kull #1 | 000001110 | | | | | | |
| Run #2 | EF155238.D | 1 | 03/24/16 | KM | 03/22/16 | OP92353 | GEF5609 |

Report of Analysis

Pesticide PPL List

950 ml

Run #2

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------|-----------------|--------|---------|-------|---|
| 309-00-2 | Aldrin | ND | 0.011 | 0.0019 | ug/l | |
| 319-84-6 | alpha-BHC | ND | 0.011 | 0.0014 | ug/l | |
| 319-85-7 | beta-BHC | ND | 0.011 | 0.0034 | ug/l | |
| 319-86-8 | delta-BHC | ND | 0.011 | 0.0017 | ug/l | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.011 | 0.00096 | ug/l | |
| 12789-03-6 | Chlordane | ND | 0.11 | 0.037 | ug/l | |
| 60-57-1 | Dieldrin | ND | 0.011 | 0.00095 | ug/l | |
| 72-54-8 | 4,4'-DDD | ND | 0.011 | 0.0013 | ug/l | |
| 72-55-9 | 4,4'-DDE | ND | 0.011 | 0.00093 | ug/l | |
| 50-29-3 | 4,4'-DDT | ND | 0.011 | 0.0027 | ug/l | |
| 72-20-8 | Endrin | ND | 0.011 | 0.0017 | ug/l | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.011 | 0.0026 | ug/l | |
| 7421-93-4 | Endrin aldehyde | ND | 0.011 | 0.0036 | ug/l | |
| 959-98-8 | Endosulfan-I | ND | 0.011 | 0.0012 | ug/l | |
| 33213-65-9 | Endosulfan-II | ND | 0.011 | 0.0018 | ug/l | |
| 76-44-8 | Heptachlor | ND | 0.011 | 0.0014 | ug/l | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.011 | 0.00082 | ug/l | |
| 72-43-5 | Methoxychlor | ND | 0.011 | 0.0038 | ug/l | |
| 8001-35-2 | Toxaphene | ND | 0.14 | 0.052 | ug/l | |
| 12674-11-2 | Aroclor 1016 | ND ^a | 0.26 | 0.049 | ug/l | |
| 11104-28-2 | Aroclor 1221 | ND ^a | 0.26 | 0.25 | ug/l | |
| 11141-16-5 | Aroclor 1232 | ND ^a | 0.26 | 0.20 | ug/l | |
| 53469-21-9 | Aroclor 1242 | ND ^a | 0.26 | 0.086 | ug/l | |
| 12672-29-6 | Aroclor 1248 | ND ^a | 0.26 | 0.081 | ug/l | |
| 11097-69-1 | Aroclor 1254 | ND ^a | 0.26 | 0.057 | ug/l | |
| 11096-82-5 | Aroclor 1260 | ND ^a | 0.26 | 0.062 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 877-09-8 | Tetrachloro-m-xylene | 86% | 86% | 21-13 | 37% | |
| 877-09-8 | Tetrachloro-m-xylene | 83% | 88% | 21-13 | 37% | |
| 2051-24-3 | Decachlorobiphenyl | 96% | 106% | 10-12 | 21% | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound

Page 1 of 2



13 of 1072

Report of Analysis

Client Sample ID:MW-5Lab Sample ID:JC16575-1Date Sampled:03/18/16Matrix:AQ - Ground WaterDate Received:03/18/16Method:EPA 608EPA 608Percent Solids:n/aProject:2002-2024 Cropsey Avenue, Brooklyn, NYNYNY

Pesticide PPL List

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 2051-24-3 | Decachlorobiphenyl | 102% | 95% | 10-121% |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

SGS

14 of 1072

ACCUTEST JC16575

4

4

Page 2 of 2

| | Report of Analysis Pa | | | | | | | | |
|--|----------------------------------|--------------------------|---|------------------|---------------------------|-----------------------|-----------------------------|--|--|
| Client San Lab Samp Matrix: Method: Project: | le ID: JC1657 AQ - C SW846 | Fround W 8151 S | ater W846 8151/351(sey Avenue, Bro | | Da Pe | I | 3/18/16 3/18/16 a | | |
| Run #1 Run #2 | File ID OA115503.D | DF 1 | Analyzed 03/28/16 | By VDT | Prep Date 03/22/16 | Prep Batch OP92340 | Analytical Batch GOA3990 | | |
| Run #1 Run #2 | Initial Volume 940 ml | Final V 10.0 m | V olume 1 | | | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|----------------------|------------------------|----------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 0.53 0.11 0.11 | 0.31 0.058 0.059 | ug/l ug/l ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| | 8 | | | 2 | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

SGS

15 of 1072 ACCUTEST JC16575

4

| Client Sample ID: | MW-5 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC16575-1 | Date Sampled: | 03/18/16 |
| Matrix: | AQ - Ground Water | Date Received: | 03/18/16 |
| | | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

Total Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|--------|------|-------|----|----------|-------------|--------------------------|--------------------------|
| Antimony | < 6.0 | 6.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Arsenic | < 3.0 | 3.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Beryllium | < 1.0 | 1.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Cadmium | < 3.0 | 3.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Chromium | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Copper | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Lead | < 3.0 | 3.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Mercury | < 0.20 | 0.20 | ug/l | 1 | 03/22/16 | 03/22/16 MS | SW846 7470A ¹ | SW846 7470A ³ |
| Nickel | 30.7 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Selenium | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Silver | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C 2 | SW846 3010A ⁴ |
| Thallium | < 2.0 | 2.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Zinc | < 20 | 20 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |

(1) Instrument QC Batch: MA38977

(2) Instrument QC Batch: MA39057

(3) Prep QC Batch: MP92703

(4) Prep QC Batch: MP92763

Page 1 of 1





| | Report of Analysis Page | | | | | | | | |
|--|--------------------------------|---------------------------|--------------------------|-----------|-------------------------|--|-----------------------------|--|--|
| Client Sar Lab Samp Matrix: Method: Project: | le ID: JC16 AQ EPA | 575-2 Ground Wa 624 | ater sey Avenue, Bro | ooklyn, N | Y | Date Sampled: Date Received: Percent Solids: | | | |
| Run #1 ^a Run #2 | File ID N253485.D | DF 1 | Analyzed 03/19/16 | By WO | Prep Date n/a | Prep Batch n/a | Analytical Batch VN10722 | | |
| Run #1 Run #2 | Purge Volum 5.0 ml | ne | | | | | | | |

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.091 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 2.1 | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 3.7 | 1.0 | 0.14 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





E = Indicates value exceeds calibration range

| Client Sample ID: Lab Sample ID: Matrix: Method: Project: | | MW-6 JC16575-2 AQ - Ground Water EPA 624 2002-2024 Cropsey Ave | enue, Brook | Date | Sampled: Received: ent Solids: | 03/18/16 03/18/16 n/a | | |
|---|--------|--|-------------|--------|--------------------------------------|-----------------------------|---|--|
| VOA PPL I | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichl | oroethene | 1.5 | 1.0 | 0.12 | ug/l | | |
| 75-69-4 | Trichl | orofluoromethane | ND | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl | chloride | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylene | es (total) | ND | 1.0 | 0.22 | ug/l | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Di | chloroethane-D4 (SUR) | 100% | | 72-1 | 25% | | |
| 2037-26-5 | Toluer | ne-D8 (SUR) | 100% | | 78-1 | 19% | | |
| 460-00-4 | 4-Bror | nofluorobenzene (SUR) | 102% | | 74-1 | 15% | | |
| 1868-53-7 | Dibroi | nofluoromethane (S) | 100% | | 79-1 | 20% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



Page 2 of 2

4.2



| | | | | | | | 1 460 1 01 2 |
|-------------------------------|-------------|-------------|-----------------|------------|-----------|-------------------|------------------|
| Client San | nple ID: MV | V-7 | | | | | |
| Lab Samp | le ID: JC1 | 6575-3 | | | Da | ate Sampled: 03 | 8/18/16 |
| Matrix: | AQ | - Ground W | ater | | Da | ate Received: 03 | 8/18/16 |
| Method: | EP | A 624 | | | Pe | ercent Solids: n/ | a |
| Project: | 200 | 2-2024 Crop | sey Avenue, Bro | ooklyn, NY | Y | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 ^a Run #2 | N253484.D | 1 | 03/19/16 | WO | n/a | n/a | VN10722 |
| | Purge Volu | me | | | | | |
| Run #1 | 5.0 ml | | | | | | |
| Run #2 | | | | | | | |

Report of Analysis

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | 1.5 | 1.0 | 0.091 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 0.34 | 1.0 | 0.12 | ug/l | J |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | 4.0 | 1.0 | 0.14 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2



E = Indicates value exceeds calibration range

| Client Samp Lab Sample Matrix: Method: Project: | | MW-7 JC16575-3 AQ - Ground Water EPA 624 2002-2024 Cropsey Av | enue, Brook | Date | Sampled: Received: ent Solids: | 03/18/16 03/18/16 n/a | | |
|---|--------|---|-------------|---------|--------------------------------------|-----------------------------|---|--|
| VOA PPL I | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichl | oroethene | 0.74 | 1.0 | 0.12 | ug/l | J | |
| 75-69-4 | Trichl | orofluoromethane | ND | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl | chloride | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylen | es (total) | ND | 1.0 | 0.22 | ug/l | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Di | ichloroethane-D4 (SUR) | 111% | | 72-1 | 25% | | |
| 2037-26-5 | Toluer | ne-D8 (SUR) | 101% | | 78-1 | 19% | | |
| 460-00-4 | | | | 74-115% | | | | |
| 1868-53-7 | Dibro | mofluoromethane (S) | 106% | | 79-1 | 20% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound

Page 2 of 2



20 of 1072

| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|--|------------------------------|----------------|----------------------------|-----------|-------------------------|--|-------------------------------|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: JC16 AQ - EPA | | c Water sey Avenue, Bro | ooklyn, N | Y | Date Sampled: Date Received: Percent Solids: | |
| Run #1 ^a Run #2 | File ID N253492.D | DF 1 | Analyzed 03/19/16 | By WO | Prep Date n/a | Prep Batcl n/a | n Analytical Batch VN10722 |
| Run #1 Run #2 | Purge Volun 5.0 ml | ie | | | | | |

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.091 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.14 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

4.4

21 of 1072

ACCUTEST JC16575

SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Dibromofluoromethane (S)

1868-53-7

| Client Samp Lab Sample Matrix: Method: Project: | | | Date | Sampled: Received: ent Solids: | 03/18/16 03/18/16 n/a | | |
|---|--|----------------------|--------------------------|--------------------------------------|------------------------------|---|--|
| VOA PPL L | ist | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 75-69-4 75-01-4 1330-20-7 | Trichloroethene Trichlorofluoromethane Vinyl chloride Xylenes (total) | ND ND ND ND | 1.0 2.0 1.0 1.0 | 0.12 0.20 0.13 0.22 | ug/l ug/l ug/l ug/l | | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 2037-26-5 460-00-4 | 1,2-Dichloroethane-D4 (SUR) Toluene-D8 (SUR) 4-Bromofluorobenzene (SUR) | 111% 100% 104% | | 72-1 78-1 74-1 | 19% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

105%

79-120%

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

Page 2 of 2



SGS

| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|--|-----------------------------|---------------------------------------|--------------------------|-----------|---------------------------|---|-----------------------------------|
| Client Sa Lab Sam Matrix: Method: Project: | A E | C16575-4 Q - Field Bl PA 625 EP | | ooklyn, N | Da Pe | ate Sampled: (ate Received: (ercent Solids: 1 | |
| Run #1 Run #2 | File ID F155737.I | DF D 1 | Analyzed 03/22/16 | By SD | Prep Date 03/21/16 | Prep Batch OP92311 | Analytical Batch EF6551 |
| Run #1 | Initial Vo 950 ml | ume Fina 1.0 1 | l Volume ml | | | | |

Run #2

ABN PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|-----------------------------|--------|-----|------|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 5.3 | 1.0 | ug/l | |
| 59-50-7 | 4-Chloro-3-methyl phenol | ND | 5.3 | 1.2 | ug/l | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 5.3 | 1.6 | ug/l | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 5.3 | 1.7 | ug/l | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 5.3 | 0.94 | ug/l | |
| 534-52-1 | 4,6-Dinitro-o-cresol | ND | 5.3 | 0.76 | ug/l | |
| 88-75-5 | 2-Nitrophenol | ND | 2.1 | 1.9 | ug/l | |
| 100-02-7 | 4-Nitrophenol | ND | 11 | 0.89 | ug/l | |
| 87-86-5 | Pentachlorophenol | ND | 5.3 | 2.0 | ug/l | |
| 108-95-2 | Phenol | ND | 2.1 | 0.52 | ug/l | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 2.1 | 1.3 | ug/l | |
| 83-32-9 | Acenaphthene | ND | 1.1 | 0.37 | ug/l | |
| 208-96-8 | Acenaphthylene | ND | 1.1 | 0.40 | ug/l | |
| 120-12-7 | Anthracene | ND | 1.1 | 0.42 | ug/l | |
| 92-87-5 | Benzidine | ND | 21 | 0.29 | ug/l | |
| 56-55-3 | Benzo(a)anthracene | ND | 1.1 | 0.37 | ug/l | |
| 50-32-8 | Benzo(a)pyrene | ND | 1.1 | 0.39 | ug/l | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 1.1 | 0.62 | ug/l | |
| 191-24-2 | Benzo(g,h,i)perylene | ND | 1.1 | 0.44 | ug/l | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 1.1 | 0.44 | ug/l | |
| 101-55-3 | 4-Bromophenyl phenyl ether | ND | 2.1 | 0.31 | ug/l | |
| 85-68-7 | Butyl benzyl phthalate | ND | 2.1 | 0.62 | ug/l | |
| 91-58-7 | 2-Chloronaphthalene | ND | 2.1 | 1.0 | ug/l | |
| 106-47-8 | 4-Chloroaniline | ND | 2.1 | 0.42 | ug/l | |
| 218-01-9 | Chrysene | ND | 1.1 | 0.26 | ug/l | |
| 111-91-1 | bis(2-Chloroethoxy)methane | ND | 2.1 | 0.69 | ug/l | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 2.1 | 0.56 | ug/l | |
| 108-60-1 | bis(2-Chloroisopropyl)ether | ND | 2.1 | 0.78 | ug/l | |
| 7005-72-3 | 4-Chlorophenyl phenyl ether | ND | 2.1 | 0.45 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 2.1 | 0.22 | ug/l | |
| 122-66-7 | 1,2-Diphenylhydrazine | ND | 2.1 | 0.48 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 2.1 | 0.17 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank N = Indicates presumptive evidence of a compound



E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

Client Sample ID:FBLab Sample ID:JC16575-4Date Sampled:03/18/16Matrix:AQ - Field Blank WaterDate Received:03/18/16Method:EPA 625EPA 625Percent Solids:n/aProject:2002-2024 Cropsey Avenue, Brooklyn, NYNYNational Solid So

ABN PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-----------|----------------------------|--------|--------|---------|-------|---|
| 106-46-7 | 1,4-Dichlorobenzene | ND | 2.1 | 0.19 | ug/l | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 2.1 | 0.91 | ug/l | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 2.1 | 0.59 | ug/l | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 2.1 | 1.3 | ug/l | |
| 53-70-3 | Dibenzo(a, h)anthracene | ND | 1.1 | 0.57 | ug/l | |
| 84-74-2 | Di-n-butyl phthalate | ND | 2.1 | 0.62 | ug/l | |
| 117-84-0 | Di-n-octyl phthalate | ND | 2.1 | 0.60 | ug/l | |
| 84-66-2 | Diethyl phthalate | ND | 2.1 | 0.41 | ug/l | |
| 131-11-3 | Dimethyl phthalate | ND | 2.1 | 0.35 | ug/l | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 2.1 | 0.70 | ug/l | |
| 206-44-0 | Fluoranthene | ND | 1.1 | 0.26 | ug/l | |
| 86-73-7 | Fluorene | ND | 1.1 | 0.47 | ug/l | |
| 118-74-1 | Hexachlorobenzene | ND | 1.1 | 0.57 | ug/l | |
| 87-68-3 | Hexachlorobutadiene | ND | 1.1 | 0.18 | ug/l | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 11 | 0.43 | ug/l | |
| 67-72-1 | Hexachloroethane | ND | 5.3 | 0.30 | ug/l | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 1.1 | 0.32 | ug/l | |
| 78-59-1 | Isophorone | ND | 2.1 | 0.62 | ug/l | |
| 91-20-3 | Naphthalene | ND | 1.1 | 0.34 | ug/l | |
| 98-95-3 | Nitrobenzene | ND | 2.1 | 0.44 | ug/l | |
| 62-75-9 | n-Nitrosodimethylamine | ND | 2.1 | 0.48 | ug/l | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 2.1 | 0.49 | ug/l | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 5.3 | 0.54 | ug/l | |
| 85-01-8 | Phenanthrene | ND | 1.1 | 0.38 | ug/l | |
| 129-00-0 | Pyrene | ND | 1.1 | 0.35 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.1 | 0.36 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 367-12-4 | 2-Fluorophenol | 40% | | 14-1 | 10% | |
| 4165-62-2 | Phenol-d5 | 28% | | 10-1 | 10% | |
| 118-79-6 | 2,4,6-Tribromophenol | 63% | | 38-145% | | |
| 4165-60-0 | Nitrobenzene-d5 | 65% | | 33-136% | | |
| 321-60-8 | 2-Fluorobiphenyl | 59% | | 35-1 | 27% | |
| 1718-51-0 | Terphenyl-d14 | 69% | | 11-1 | 39% | |

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

SGS

4.4 **4**

Page 2 of 2

24 of 1072 ACCUTEST JC16575

5.0 ml

| Lab Samı Matrix: Method: | AQ - Fi EPA 60 | JC16575-4 AQ - Field Blank Water EPA 608 EPA 608 | | | | L | | |
|---|---------------------------------|--|----------|----|-----------|------------|------------------|--|
| Project: 2002-2024 Cropsey Avenue, Brooklyn, NY | | | | Y | | | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | |
| Run #1 | 6G33313.D | 1 | 03/24/16 | BP | 03/22/16 | OP92352 | G6G969 | |
| Run #2 | EF155239.D | 1 | 03/24/16 | KM | 03/22/16 | OP92353 | GEF5609 | |
| Run #1 | Initial Volume 970 ml | Final V 5.0 ml | Volume | | | | | |

Report of Analysis

Pesticide PPL List

970 ml

Run #2

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------|-----------------|--------|---------|-------|---|
| 309-00-2 | Aldrin | ND | 0.010 | 0.0017 | ug/l | |
| 319-84-6 | alpha-BHC | ND | 0.010 | 0.0013 | ug/l | |
| 319-85-7 | beta-BHC | ND | 0.010 | 0.0032 | ug/l | |
| 319-86-8 | delta-BHC | ND | 0.010 | 0.0016 | ug/l | |
| 58-89-9 | gamma-BHC (Lindane) | ND | 0.010 | 0.00089 | ug/l | |
| 12789-03-6 | Chlordane | ND | 0.10 | 0.034 | ug/l | |
| 60-57-1 | Dieldrin | ND | 0.010 | 0.00088 | ug/l | |
| 72-54-8 | 4,4'-DDD | ND | 0.010 | 0.0012 | ug/l | |
| 72-55-9 | 4,4'-DDE | ND | 0.010 | 0.00086 | ug/l | |
| 50-29-3 | 4,4'-DDT | ND | 0.010 | 0.0025 | ug/l | |
| 72-20-8 | Endrin | ND | 0.010 | 0.0016 | ug/l | |
| 1031-07-8 | Endosulfan sulfate | ND | 0.010 | 0.0024 | ug/l | |
| 7421-93-4 | Endrin aldehyde | ND | 0.010 | 0.0033 | ug/l | |
| 959-98-8 | Endosulfan-I | ND | 0.010 | 0.0011 | ug/l | |
| 33213-65-9 | Endosulfan-II | ND | 0.010 | 0.0017 | ug/l | |
| 76-44-8 | Heptachlor | ND | 0.010 | 0.0013 | ug/l | |
| 1024-57-3 | Heptachlor epoxide | ND | 0.010 | 0.00076 | ug/l | |
| 72-43-5 | Methoxychlor | ND | 0.010 | 0.0035 | ug/l | |
| 8001-35-2 | Toxaphene | ND | 0.13 | 0.049 | ug/l | |
| 12674-11-2 | Aroclor 1016 | ND a | 0.26 | 0.048 | ug/l | |
| 11104-28-2 | Aroclor 1221 | ND a | 0.26 | 0.24 | ug/l | |
| 11141-16-5 | Aroclor 1232 | ND a | 0.26 | 0.20 | ug/l | |
| 53469-21-9 | Aroclor 1242 | ND a | 0.26 | 0.085 | ug/l | |
| 12672-29-6 | Aroclor 1248 | ND ^a | 0.26 | 0.079 | ug/l | |
| 11097-69-1 | Aroclor 1254 | ND ^a | 0.26 | 0.056 | ug/l | |
| 11096-82-5 | Aroclor 1260 | ND ^a | 0.26 | 0.060 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 877-09-8 | Tetrachloro-m-xylene | 83% | 95% | 21-13 | 37% | |
| 877-09-8 | Tetrachloro-m-xylene | 79% | 96% | 21-13 | 37% | |
| 2051-24-3 | Decachlorobiphenyl | 53% | 69% | 10-12 | 21% | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



Page 1 of 2



Report of Analysis

Client Sample ID:FBLab Sample ID:JC16575-4Date Sampled:03/18/16Matrix:AQ - Field Blank WaterDate Received:03/18/16Method:EPA 608EPA 608Percent Solids:n/aProject:2002-2024 Cropsey Avenue, Brooklyn, NYProject:Client Solids:n/a

Pesticide PPL List

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 2051-24-3 | Decachlorobiphenyl | 58% | 63% | 10-121% |

(a) Result is from Run# 2

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





Page 2 of 2



26 of 1072 ACCUTEST JC16575

| | | | Repo | ort of A | nalysis | | Page 1 of |
|--------------------------------|--------------------------------|----------------------|---|------------------|---------------------------|-----------------------|-----------------------------|
| Client Sar Lab Samj | ple ID: JC1 | 5575-4 | L W/ A | | | ate Sampled: 03 | 3/18/16 |
| Matrix: Method: Project: | swe | | k water W846 8151/351 sey Avenue, Bro | | Pe | ercent Solids: n/ | |
| Run #1 Run #2 | File ID OA115476.E | DF 0 1 | Analyzed 03/26/16 | By VDT | Prep Date 03/22/16 | Prep Batch OP92340 | Analytical Batch GOA3989 |
| Run #1 Run #2 | Initial Volun 800 ml | ne Final V 10.0 m | V olume 1 | | | | |

Herbicide List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|-------------------------------|---------------------------------------|----------------|----------------------|------------------------|----------------------|---|
| 94-75-7 93-72-1 93-76-5 | 2,4-D 2,4,5-TP (Silvex) 2,4,5-T | ND ND ND | 0.63 0.13 0.13 | 0.36 0.069 0.069 | ug/l ug/l ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 19719-28-9 19719-28-9 | 2,4-DCAA 2,4-DCAA | 85% 67% | | 39-1: 39-1: | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



27 of 1072 ACCUTEST JC16575



| Client Sample ID: | FB | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC16575-4 | Date Sampled: | 03/18/16 |
| Matrix: | AQ - Field Blank Water | Date Received: | 03/18/16 |
| | | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

Report of Analysis

Total Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|--------|------|-------|----|----------|-------------|--------------------------|--------------------------|
| Antimony | < 6.0 | 6.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Arsenic | < 3.0 | 3.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Beryllium | < 1.0 | 1.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Cadmium | < 3.0 | 3.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Chromium | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Copper | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Lead | < 3.0 | 3.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Mercury | 2.5 | 0.20 | ug/l | 1 | 03/22/16 | 03/22/16 MS | SW846 7470A ¹ | SW846 7470A ³ |
| Nickel | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Selenium | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Silver | < 10 | 10 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Thallium | < 2.0 | 2.0 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C ² | SW846 3010A ⁴ |
| Zinc | < 20 | 20 | ug/l | 1 | 03/24/16 | 03/30/16 GT | SW846 6010C 2 | SW846 3010A ⁴ |

(1) Instrument QC Batch: MA38977

(2) Instrument QC Batch: MA39057

(3) Prep QC Batch: MP92703

(4) Prep QC Batch: MP92763



4.4



| Report of Analysis Pa | | | | | | | | | |
|--|-----------------------------|---|--------------------------|-----------|-------------------------|--|------------------------------------|--|--|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: JC1 AQ EPA | 6575-5 - Trip Blank A 624 2-2024 Crops | Water sey Avenue, Bro | ooklyn, N | Y | Date Sampled: (Date Received: (Percent Solids:) | | | |
| Run #1 ^a Run #2 | File ID N253493.D | DF 1 | Analyzed 03/19/16 | By WO | Prep Date n/a | Prep Batch n/a | Analytical Batch VN10722 | | |
| Run #1 Run #2 | Purge Volu 5.0 ml | me | | | | | | | |

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|-------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.6 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 2.6 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.10 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.10 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.17 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.57 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.096 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.093 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.21 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.50 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.091 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.11 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.15 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.29 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.12 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.090 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.16 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.12 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.11 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.12 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.15 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.22 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.22 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.12 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.14 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.086 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.13 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



29 of 1072

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

1868-53-7

| Client Samp Lab Sample Matrix: Method: Project: | | TB JC16575-5 AQ - Trip Blank Water EPA 624 2002-2024 Cropsey Ave | | lyn, NY | | Date | Sampled: Received: ent Solids: | 03/18/16 03/18/16 n/a |
|---|--------|--|--------|---------|---------------|-------|--------------------------------------|-----------------------------|
| VOA PPL I | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | |
| 79-01-6 | Trichl | oroethene | ND | 1.0 | 0.12 | ug/l | | |
| 75-69-4 | Trichl | orofluoromethane | ND | 2.0 | 0.20 | ug/l | | |
| 75-01-4 | Vinyl | chloride | ND | 1.0 | 0.13 | ug/l | | |
| 1330-20-7 | Xylene | es (total) | ND | 1.0 | 1.0 0.22 ug/l | | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 | 1,2-Di | chloroethane-D4 (SUR) | 110% | | | | | |
| 2037-26-5 | Toluer | ne-D8 (SUR) | 102% | | 78-1 | 19% | | |
| 460-00-4 | 4-Bror | nofluorobenzene (SUR) | 103% | | 74-1 | 15% | | |

Report of Analysis

(a) (pH= 5) Sample is not acid preserved per method/client criteria. Sample analyzed within 3 days holding time as required for acrolein and acrylonitrile. Other compounds within 7 days as required by the method.

79-120%

108%

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

Page 2 of 2



30 of 1072



Section 4

Sample Results

Report of Analysis





| Client San Lab Samp Matrix: Method: Project: | le ID: JC A(EP | IP BLANK 17397-1 2 - Trip Blank A 624)2-2024 Crops | Water sey Avenue, Bro | ooklyn, NY | Da | ate Sampled: 03 ate Received: 03 ercent Solids: n/ | | |
|--|-----------------------------|---|--------------------------|------------|-------------------------|--|-----------------------------|--|
| Run #1 ^a Run #2 | File ID N253844.D | DF 1 | Analyzed 04/04/16 | By WO | Prep Date n/a | Prep Batch n/a | Analytical Batch VN10735 | |
| Run #1 Run #2 | Purge Volu 5.0 ml | me | | | | | | |

Report of Analysis

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.3 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 1.2 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.11 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.14 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.15 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.20 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.19 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.12 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.19 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.73 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.11 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.22 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.15 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.15 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.13 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.54 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.16 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.22 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.30 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.17 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.16 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.17 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.16 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.18 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.17 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.12 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.17 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.18 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $J = \ Indicates \ an \ estimated \ value$

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



4.1

Page 1 of 2

7 of 180

ACCUTEST JC17397

SGS

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

1868-53-7

| Client Samp Lab Sample Matrix: Method: Project: | | TRIP BLANK JC17397-1 AQ - Trip Blank Water EPA 624 2002-2024 Cropsey Ave | | lyn, NY | | Date Sampled:03/31/16Date Received:03/31/16Percent Solids:n/a | | | |
|---|--------|--|----------|------------|----------------------|---|---|--|--|
| VOA PPL I | List | | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDL | Units | Q | | |
| 79-01-6 | | proethene | ND | 1.0 | 0.16 | ug/l | | | |
| 75-69-4 75-01-4 | | orofluoromethane chloride | ND ND | 2.0 1.0 | 0.55 0.19 | ug/l ug/l | | | |
| 1330-20-7 | Xylene | es (total) | ND | 1.0 | 0.24 | ug/l | | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | Lim | its | | | |
| 17060-07-0 2037-26-5 460-00-4 | Toluer | chloroethane-D4 (SUR) ne-D8 (SUR) nofluorobenzene (SUR) | 99% | | 72-1 78-1 74-1 | 19% | | | |

Report of Analysis

(a) Results reported from the HCl preserved sample. This reported result can only be used for screening purposes for acrolein and acrylonitrile.

79-120%

100%

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2



| Report of Analysis Page | | | | | | | | | | |
|--|------------------------------|----------------|--------------------------|----------|----------------------|--|-------------------------------|--|--|--|
| Client Sar Lab Samp Matrix: Method: Project: | ole ID: JC173 AQ - EPA | | | | | Date Sampled: Date Received: Percent Solids: | | | | |
| Run #1 ^a Run #2 | File ID N253845.D | DF 1 | Analyzed 04/04/16 | By WO | Prep Date n/a | Prep Batc n/a | h Analytical Batch VN10735 | | | |
| Run #1 Run #2 | Purge Volum 5.0 ml | e | | | | | | | | |

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|--------|-----|------|-------|---|
| 107-02-8 | Acrolein | ND | 10 | 1.3 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 10 | 1.2 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | 0.11 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.14 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.15 | ug/l | |
| 74-83-9 | Bromomethane | ND | 1.0 | 0.20 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.19 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.12 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.19 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 5.0 | 0.73 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.11 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.22 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.15 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.15 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.13 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.54 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.16 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.22 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.30 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.14 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.17 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.16 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.17 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.11 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 1.0 | 0.16 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.18 | ug/l | |
| 127-18-4 | Tetrachloroethene | 0.83 | 1.0 | 0.17 | ug/l | J |
| 108-88-3 | Toluene | ND | 1.0 | 0.12 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.17 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.18 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



9 of 180

ACCUTEST JC17397

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

1868-53-7

| Client Samp Lab Sample Matrix: Method: Project: | ID: JC17397-2 AQ - Field I EPA 624 | Blank Water Cropsey Avenue, Broo | oklyn, NY | | Date | Sampled: Received: ent Solids: | 03/31/16 03/31/16 n/a |
|---|---|-------------------------------------|--------------------------|------------------------------|------------------------------|--------------------------------------|-----------------------------|
| VOA PPL L | ist | | | | | | |
| CAS No. | Compound | Result | RL | MDL | Units | Q | |
| 79-01-6 75-69-4 75-01-4 1330-20-7 | Trichloroethene Trichlorofluorometh Vinyl chloride Xylenes (total) | ND ND ND ND | 1.0 2.0 1.0 1.0 | 0.16 0.55 0.19 0.24 | ug/l ug/l ug/l ug/l | | |
| CAS No. | Surrogate Recover | ies Run# 1 | Run# 2 | Lim | its | | |
| 17060-07-0 2037-26-5 460-00-4 | 1,2-Dichloroethane- Toluene-D8 (SUR) 4-Bromofluorobenz | 100% | | 72-1 78-1 74-1 | 19% | | |

Report of Analysis

(a) Results reported from the HCl preserved sample. This reported result can only be used for screening purposes for acrolein and acrylonitrile.

104%

79-120%

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2



10 of 180

ACCUTEST JC17397

| Lab Samp Matrix: Method: Project: | AQ - EPA | JC17397-3 AQ - Ground Water EPA 624 2002-2024 Cropsey Avenue, Brooklyn, NY | | | | Date Sampled:03/31/16Date Received:03/31/16Percent Solids:n/a | | | |
|--|-------------|---|----------|----|-----------|---|------------------|--|--|
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch | | |
| Run #1 ^a | N253846.D | 10 | 04/05/16 | WO | n/a | n/a | VN10735 | | |
| Run #2 ^a | N253847.D | 100 | 04/05/16 | WO | n/a | n/a | VN10735 | | |
| | | | | | | | | | |

Report of Analysis

Run #1 5.0 ml

Run #2 5.0 ml

VOA PPL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|---------------------------|-------------------|-----|-----|-------|---|
| 107-02-8 | Acrolein | ND | 100 | 13 | ug/l | |
| 107-13-1 | Acrylonitrile | ND | 100 | 12 | ug/l | |
| 71-43-2 | Benzene | ND | 10 | 1.1 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 10 | 1.4 | ug/l | |
| 75-25-2 | Bromoform | ND | 10 | 1.5 | ug/l | |
| 74-83-9 | Bromomethane | ND | 10 | 2.0 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 10 | 1.9 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 10 | 1.2 | ug/l | |
| 75-00-3 | Chloroethane | ND | 10 | 1.9 | ug/l | |
| 110-75-8 | 2-Chloroethyl vinyl ether | ND | 50 | 7.3 | ug/l | |
| 67-66-3 | Chloroform | ND | 10 | 1.1 | ug/l | |
| 74-87-3 | Chloromethane | ND | 10 | 2.2 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 10 | 2.2 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 10 | 1.5 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 10 | 1.5 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 10 | 1.3 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 20 | 5.4 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 10 | 1.6 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 10 | 2.1 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 10 | 2.2 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 924 | 10 | 3.0 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | 8.0 | 10 | 1.4 | ug/l | J |
| 78-87-5 | 1,2-Dichloropropane | ND | 10 | 1.7 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 10 | 1.6 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 10 | 1.7 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 10 | 1.1 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 10 | 1.6 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 10 | 1.8 | ug/l | |
| 127-18-4 | Tetrachloroethene | 7270 ^b | 100 | 17 | ug/l | |
| 108-88-3 | Toluene | ND | 10 | 1.2 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 10 | 1.7 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 10 | 1.8 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





SGS

E = Indicates value exceeds calibration range

Dibromofluoromethane (S)

| Client Samp Lab Sample Matrix: Method: Project: | | MW-5 JC17397-3 AQ - Ground Water EPA 624 2002-2024 Cropsey Ave | enue, Brook | Date | Sampled: Received: ent Solids: | 03/31/16 03/31/16 n/a | | |
|---|--------|--|-------------|--------|--------------------------------------|-----------------------------|---|--|
| VOA PPL I | List | | | | | | | |
| CAS No. | Comp | ound | Result | RL | MDI | L Units | Q | |
| 79-01-6 | Trichl | oroethene | 181 | 10 | 1.6 | ug/l | | |
| 75-69-4 | Trichl | orofluoromethane | ND | 20 | 5.5 | ug/l | | |
| 75-01-4 | Vinyl | chloride | ND | 10 | 1.9 | ug/l | | |
| 1330-20-7 | Xylen | es (total) | ND | 10 | 2.4 | ug/l | | |
| CAS No. | Surro | gate Recoveries | Run# 1 | Run# 2 | L | imits | | |
| 17060-07-0 | 1,2-Di | chloroethane-D4 (SUR) | 91% | 97% | 7 | 2-125% | | |
| 2037-26-5 | Toluer | ne-D8 (SUR) | 97% | 99% | 7 | 8-119% | | |
| 460-00-4 | 4-Broi | nofluorobenzene (SUR) | 97% | 98% | 74 | 4-115% | | |

Report of Analysis

(a) Results reported from the HCl preserved sample. This reported result can only be used for screening purposes for acrolein and acrylonitrile.

99%

79-120%

96%

(b) Result is from Run# 2

1868-53-7

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



12 of 180

JC17397





Section 4

Sample Results

Report of Analysis





| Report of Analysis Page | | | | | | | | | | |
|---|------------------------------|------------------------------------|--------------------------|-----------------|----------------------|--|-------------------------------|--|--|--|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC2 AQ SW | 3746-1 - Ground Wa 846 8260C | ater sey Avenue, Bro | ooklyn, N | Y | Date Sampled: Date Received: Percent Solids: | 0.11 0.01 - 0 | | | |
| Run #1 Run #2 | File ID 3B128390.D | DF 1 | Analyzed 07/11/16 | By HA | Prep Date n/a | Prep Batcl n/a | n Analytical Batch V3B5741 | | | |
| Run #1 Run #2 | Purge Volu 5.0 ml | me | | | | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | 0.78 | 1.0 | 0.23 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 1.3 | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



JC23746

SGS

E = Indicates value exceeds calibration range

Report of Analysis

| Client Sample ID: | MW-5I | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC23746-1 | Date Sampled: | 07/08/16 |
| Matrix: | AQ - Ground Water | Date Received: | 07/08/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | 18.7 | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | 0.47 | 1.0 | 0.22 | ug/l | J |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | 2.1 | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 99% | | 76-1 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 100% | | 73-1 | 22% | |
| 2037-26-5 | Toluene-D8 | 100% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 99% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|---|-----------------------------|--|--------------------------|-----------|-------------------------|--|-----------------------------|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC A(SV | W-8 23746-2 Q - Ground Wa V846 8260C 02-2024 Crops | ater sey Avenue, Bro | ooklyn, N | | Date Sampled: Date Received: Percent Solids: | 011 001 20 |
| Run #1 | File ID 3B128391. | DF D 1 | Analyzed 07/11/16 | By HA | Prep Date n/a | Prep Batch n/a | Analytical Batch V3B5741 |
| Run #2 Run #1 Run #2 | Purge Vol 5.0 ml | ume | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.23 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



9 of 194

JC23746

4.2

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Client Sample ID: | MW-8 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC23746-2 | Date Sampled: | 07/08/16 |
| Matrix: | AQ - Ground Water | Date Received: | 07/08/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| - | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | 0.52 | 1.0 | 0.23 | ug/l | J |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.22 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 99% | | 76-1 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 103% | | 73-1 | 22% | |
| 2037-26-5 | Toluene-D8 | 100% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 99% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| | | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|---|---------------------------|-------|----------------|--------------------------|-----------|-------------------------|--|-------------------------------|
| Client San Lab Sam Matrix: Method: Project: | ple ID: | SW846 | Fround Wa | nter sey Avenue, Bro | ooklyn, N | | Date Sampled: Date Received: Percent Solids: | 0.1.001-0 |
| Run #1 Run #2 | File ID 3B12839 | 2.D | DF 1 | Analyzed 07/11/16 | By HA | Prep Date n/a | Prep Batc n/a | h Analytical Batch V3B5741 |
| Run #1 | Purge V 5.0 ml | olume | | | | | | |

Run #2

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | 0.66 | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.23 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 1.1 | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



SGS

E = Indicates value exceeds calibration range

Report of Analysis

| Client Sample ID: | MW-9 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC23746-3 | Date Sampled: | 07/08/16 |
| Matrix: | AQ - Ground Water | Date Received: | 07/08/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|---------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | 7.5 | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.22 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | 0.87 | 1.0 | 0.26 | ug/l | J |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 100% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 103% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 102% | 84-119% | | | |
| 460-00-4 | 4-Bromofluorobenzene | 100% | 78-117% | | | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





12 of 194

JC23746

| Report of Analysis Pa | | | | | | | |
|---|------------------------------|---|--------------------------|-----------------|-------------------------|--|-----------------------------|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC A(SW | W-10 23746-4 2 - Ground Wa V846 8260C 02-2024 Crops | ater sey Avenue, Bro | ooklyn, N | | Date Sampled: Date Received: Percent Solids: | 011 001 20 |
| Run #1 Run #2 | File ID 3B128491.7 | DF D 1 | Analyzed 07/14/16 | By HA | Prep Date n/a | Prep Batch n/a | Analytical Batch V3B5745 |
| Run #1 Run #2 | Purge Volu 5.0 ml | ıme | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | 0.33 | 1.0 | 0.23 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 2.9 | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



SGS

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Client Sample ID: | MW-10 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC23746-4 | Date Sampled: | 07/08/16 |
| Matrix: | AQ - Ground Water | Date Received: | 07/08/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| 0 | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | 17.6 | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.22 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | 3.4 | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 101% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 106% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 102% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 99% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|---|------------------------------|---|--------------------------|-----------------|-------------------------|--|-----------------------------|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC: AQ SW | V-100 23746-5 2 - Ground Wa 846 8260C 92-2024 Crops | ater sey Avenue, Bro | ooklyn, N | | Date Sampled: Date Received: Percent Solids: | 011 001 20 |
| Run #1 Run #2 | File ID 3B128492.1 | DF D 1 | Analyzed 07/14/16 | By HA | Prep Date n/a | Prep Batch n/a | Analytical Batch V3B5745 |
| Run #1 Run #2 | Purge Volu 5.0 ml | me | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | 0.72 | 1.0 | 0.23 | ug/l | J |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | 1.4 | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



SGS

E = Indicates value exceeds calibration range

Report of Analysis

| Client Sample ID: | MW-100 | | |
|-------------------|--|-----------------|----------|
| Lab Sample ID: | JC23746-5 | Date Sampled: | 07/08/16 |
| Matrix: | AQ - Ground Water | Date Received: | 07/08/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |
| | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | 18.4 | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | 0.49 | 1.0 | 0.22 | ug/l | J |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | 2.1 | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Lim | its | |
| 1868-53-7 | Dibromofluoromethane | 101% | | 76-1 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 106% | | 73-1 | 22% | |
| 2037-26-5 | Toluene-D8 | 103% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 96% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2

SGS

| | | | Repo | ort of A | Analysis | | Page 1 of 2 |
|---|------------------------------|--|----------------------------|-----------------|-------------------------|--|------------------------------------|
| Client San Lab Sam Matrix: Method: Project: | ple ID: JC2 AQ SW | 3746-6 - Field Blank 346 8260C | s Water sey Avenue, Bro | ooklyn, N | | Date Sampled: Date Received: Percent Solids: | |
| Run #1 Run #2 | File ID 3B128396.D | DF 1 | Analyzed 07/11/16 | By HA | Prep Date n/a | Prep Batch n/a | Analytical Batch V3B5741 |
| Run #1 Run #2 | Purge Volu 5.0 ml | ne | | | | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.23 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



17 of 194

ACCUTEST JC23746

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Report of Analysis

| Client Sample ID: | FIELD | | |
|--------------------------|--|-----------------|----------|
| Lab Sample ID: | JC23746-6 | Date Sampled: | 07/08/16 |
| Matrix: | AQ - Field Blank Water | Date Received: | 07/08/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.22 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | its | |
| 1868-53-7 | Dibromofluoromethane | 99% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 101% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 100% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 98% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2



18 of 194

ACCUTEST JC23746

| Client Sam Lab Samp Matrix: Method: | le ID: JC23 AQ - | 746-7 Trip Blank 46 8260C | Water | | Da | ate Sampled: 07 ate Received: 07 ercent Solids: n/ | |
|--|---------------------|---------------------------------|-----------------|------------|-----------|--|------------------|
| Project: | 2002 | -2024 Crop | sey Avenue, Bro | ooklyn, NY | , | | |
| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
| Run #1 | 3B128397.D | 1 | 07/11/16 | HA | n/a | n/a | V3B5741 |
| Run #2 | | | | | | | |
| | Purge Volun | ne | | | | | |
| Run #1 | 5.0 ml | | | | | | |
| | 2.0 ml | | | | | | |

Report of Analysis

Run #2

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|-----------------------------|--------|------|------|-------|---|
| 67-64-1 | Acetone | ND | 10 | 3.8 | ug/l | |
| 71-43-2 | Benzene | ND | 0.50 | 0.14 | ug/l | |
| 74-97-5 | Bromochloromethane | ND | 1.0 | 0.46 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | 0.55 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | 0.34 | ug/l | |
| 74-83-9 | Bromomethane | ND | 2.0 | 0.46 | ug/l | |
| 78-93-3 | 2-Butanone (MEK) | ND | 10 | 1.9 | ug/l | |
| 75-15-0 | Carbon disulfide | ND | 2.0 | 0.33 | ug/l | |
| 56-23-5 | Carbon tetrachloride | ND | 1.0 | 0.54 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | 0.17 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | 0.44 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | 0.23 | ug/l | |
| 74-87-3 | Chloromethane | ND | 1.0 | 0.96 | ug/l | |
| 110-82-7 | Cyclohexane | ND | 5.0 | 0.73 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 2.0 | 0.69 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | 0.23 | ug/l | |
| 106-93-4 | 1,2-Dibromoethane | ND | 1.0 | 0.22 | ug/l | |
| 95-50-1 | 1,2-Dichlorobenzene | ND | 1.0 | 0.23 | ug/l | |
| 541-73-1 | 1,3-Dichlorobenzene | ND | 1.0 | 0.19 | ug/l | |
| 106-46-7 | 1,4-Dichlorobenzene | ND | 1.0 | 0.21 | ug/l | |
| 75-71-8 | Dichlorodifluoromethane | ND | 2.0 | 0.70 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | 0.21 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | 0.39 | ug/l | |
| 75-35-4 | 1,1-Dichloroethene | ND | 1.0 | 0.20 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethene | ND | 1.0 | 0.31 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethene | ND | 1.0 | 0.36 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | 0.33 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | 0.19 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | 0.26 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | 0.20 | ug/l | |
| 76-13-1 | Freon 113 | ND | 5.0 | 1.2 | ug/l | |
| 591-78-6 | 2-Hexanone | ND | 5.0 | 1.5 | ug/l | |

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$

N = Indicates presumptive evidence of a compound



Page 1 of 2



E = Indicates value exceeds calibration range

Report of Analysis

| Client Sample ID: | TRIP | | |
|-------------------|--|-----------------------|----------|
| Lab Sample ID: | JC23746-7 | Date Sampled: | 07/08/16 |
| Matrix: | AQ - Trip Blank Water | Date Received: | 07/08/16 |
| Method: | SW846 8260C | Percent Solids: | n/a |
| Project: | 2002-2024 Cropsey Avenue, Brooklyn, NY | | |

VOA TCL List

| CAS No. | Compound | Result | RL | MDL | Units | Q |
|------------|----------------------------|--------|--------|-------|-------|---|
| 98-82-8 | Isopropylbenzene | ND | 1.0 | 0.16 | ug/l | |
| 79-20-9 | Methyl Acetate | ND | 5.0 | 1.5 | ug/l | |
| 108-87-2 | Methylcyclohexane | ND | 5.0 | 0.78 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | 0.34 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone(MIBK) | ND | 5.0 | 1.2 | ug/l | |
| 75-09-2 | Methylene chloride | ND | 2.0 | 0.35 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | 0.27 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 1.0 | 0.39 | ug/l | |
| 127-18-4 | Tetrachloroethene | ND | 1.0 | 0.23 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | 0.23 | ug/l | |
| 87-61-6 | 1,2,3-Trichlorobenzene | ND | 1.0 | 0.20 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 1.0 | 0.25 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | 0.22 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | 0.28 | ug/l | |
| 79-01-6 | Trichloroethene | ND | 1.0 | 0.26 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 2.0 | 0.58 | ug/l | |
| 75-01-4 | Vinyl chloride | ND | 1.0 | 0.33 | ug/l | |
| | m,p-Xylene | ND | 1.0 | 0.42 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | 0.21 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 1.0 | 0.21 | ug/l | |
| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limi | ts | |
| 1868-53-7 | Dibromofluoromethane | 100% | | 76-12 | 20% | |
| 17060-07-0 | 1,2-Dichloroethane-D4 | 101% | | 73-12 | 22% | |
| 2037-26-5 | Toluene-D8 | 102% | | 84-1 | 19% | |
| 460-00-4 | 4-Bromofluorobenzene | 98% | | 78-1 | 17% | |

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Page 2 of 2

JC23746

Appendix H

Data Usability Summary Report (DUSR)

APEX COMPANIES, LLC

RI Appendix H Data Usability Summary Report



Cropsey Site 2002-2024 Cropsey Avenue Brooklyn, NY

PREPARED FOR:

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March 9, 2018

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Table of Contents (Appendix H Page Numbers Preceded by H-)

| Sco | pe o | of Work | 1 |
|-----|------|--|----|
| Sco | pe o | of Work | 1 |
| 1.0 | | Analytical Laboratory | 1 |
| | 1.1 | TNI Accreditation | 1 |
| | 1.2 | Laboratory Data Packages | 1 |
| | 1.3 | Laboratory Analytical Methods | 1 |
| | | 1.3.1 Volatile Organic Analysis | 1 |
| | | 1.3.2 Semi-Volatile Organic Analysis | 2 |
| | | 1.3.3 Pesticides, PCBs, and Herbicides | 2 |
| | | 1.3.4 Metals and Mercury | 2 |
| | 1.4 | Data Qualification | 2 |
| 2.0 | | Data Validation | 3 |
| | 2.1 | SDG JC7512 | 3 |
| | | 2.1.1 VOCs | 3 |
| | | 2.1.2 SVOCs | 5 |
| | | 2.1.3 Pesticides | 7 |
| | | 2.1.4 PCBs | 8 |
| | | 2.1.5 Herbicides | 8 |
| | | 2.1.6 Metals | 9 |
| | | 2.1.7 Mercury | 10 |
| | | 2.1.8 Percent Moisture | 10 |
| | 2.2 | SDG JC7519 | 10 |
| | 2.3 | SDG JC7667 | 12 |
| | 2.4 | SDG JC12861 | 13 |
| | | 2.4.1 VOCs | 14 |
| | | 2.4.2 Percent Moisture | 25 |
| | 2.5 | SDG JC14061 | 25 |
| | | 2.5.1 VOCs | 26 |
| | | 2.5.2 SVOCs | 27 |
| | | 2.5.3 Pesticides | 29 |
| | | 2.5.4 PCBS | 30 |
| | | 2.5.5 Herbicides | 30 |

| | 2.5.6 | Metals | |
|------|---------|------------------|----|
| | 2.5.7 | Mercury | |
| 2.6 | SDG J | C15505 | |
| 2.7 | SDG J | C15742 | |
| | 2.7.1 | VOCs | |
| | 2.7.2 | Percent Moisture | |
| 2.8 | SDG J | C15931 | |
| 2.9 | SDG J | C16088 | |
| | 2.9.1 | VOCs | |
| | 2.9.2 | Percent Moisture | |
| 2.10 |) SDG J | C16571 | 41 |
| 2.1 | I SDG J | C16575 | 42 |
| | 2.11.1 | VOCs | |
| | 2.11.2 | SVOCs | |
| | 2.11.3 | Pesticides | |
| | 2.11.4 | PCBS | |
| | 2.11.5 | Herbicides | |
| | 2.11.6 | Metals | |
| | 2.11.7 | Mercury | |
| 2.12 | 2 SDG J | C17397 | |
| | 2.12.1 | VOCs | |
| 2.13 | B SDG J | C22892 | 51 |
| | 2.13.1 | VOCs | |
| | 2.13.2 | Percent Moisture | |
| 2.14 | 1 SDG J | C23746 | |
| | 2.14.1 | VOCs | |
| | | | |

Data Usability Summary Report BCP Site No. C224169 2002-2024 Cropsey Avenue Brooklyn, NY

Scope of Work

Environmental samples were collected from the Cropsey site in Brooklyn, New York. These samples consisted of groundwater, soil, soil vapor, and ambient air which were collected between October 29, 2015 and July 8, 2016. Analytical results were validated and usability was determined using the following guidelines:

- NYSDEC Analytical Services Protocol (ASP);
- USEPA Contract Laboratory Program (CLP) National Functional Guidelines for Organic and Inorganic Data Validation, and;
- USEPA analytical methods.

1.0 Analytical Laboratory

Samples collected were analyzed by SGS Accutest, Inc. of Dayton, New Jersey. SGS Accutest holds accreditation under The NELAC Institute (TNI) lab code TNI01283.

1.1 TNI Accreditation

SGS Accutest, Inc. was accredited for all testing performed except for 1,2-dichlorobenzene, 1,3-dichlorobenzene and 1,4-dichlorobenzene by EPA 625 from laboratory report JC16575. These same analytes were analyzed by EPA 624 in JC16575, which was included in SGS Accutest's accreditation at the time, so reportable data completeness was not affected.

1.2 Laboratory Data Packages

Laboratory data packages were complete and included cover pages, chain of custodies, sample log-in information, and case narratives. Raw data was included in data packages and was used in the validation process.

1.3 Laboratory Analytical Methods

Analyses of soil and groundwater samples collected consisted of volatile organic compounds (VOC), semi-volatile organic compounds (SVOC), chlorinated pesticides, chlorinated herbicides, metals, mercury, polychlorinated biphenyls (PCB) and percent moisture (soil only). Soil samples were collected using EnCore or similar sampling devices for volatile organic compounds (VOC) under EPA 5035. Soil vapor and ambient air samples were analyzed for VOCs. Data validation for these analytical methods are included in Section 2.

1.3.1 Volatile Organic Analysis

Soil samples were prepared by EPA 5035 and analyzed by EPA 8260C. Groundwater samples were generally analyzed by EPA 624, apart from SDG JC23746, which was analyzed by EPA 8260C. Both methods are acceptable for groundwater analysis. Air samples were analyzed by TO-15. Data validation included review of requirements from EPA

methods and recommendations from the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 (CLP Guidance). Groundwater and air samples were reviewed under the "Trace Volatiles" CLP Guidance; while soil VOCs were reviewed under the "Low/Medium Volatiles" CLP Guidance.

1.3.2 Semi-Volatile Organic Analysis

Soil samples were analyzed by EPA 8270D and groundwater samples were analyzed by EPA 625. Data was validated using EPA methods 8270D and EPA 625 and the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 (CLP Guidance).

1.3.3 Pesticides, PCBs, and Herbicides

EPA methods 8081B, 8082A and 8151 were used to accurately analyze soil samples. Groundwater samples were analyzed by EPA 608 and 8151. CLP Guidance and EPA methods 8081, 8082, 608 and 8151 were used to validate data.

1.3.4 Metals and Mercury

Soil an groundwater samples were analyzed by EPA 6010C. Mercury was analyzed by 7471B and 7470, depending on the matrix. These results were validated using EPA *National Functional Guidelines for Superfund Inorganic Methods Review*, January 2017 (CLP Guidance).

1.4 Data Qualification

Laboratory data may be qualified with the following flags:

- J The analyte result is an estimated value.
- **U** The analyte was analyzed for but was not detected over the laboratory detection limit.
- **UJ** The analyte was analyzed for but was not detected over the laboratory detection limit. Due to findings with data quality, the laboratory detection limit may be inaccurate and is an estimated value.
- **R** The analyte result is unusable due to data quality deficiencies.

2.0 Data Validation

The following subsections include data validation for individual Sample Delivery Groups (SDG). Each sample within the SDG was reviewed based on EPA method guidelines, CLP Guidance's and NYSDEC's ASP. Summaries of data validation findings and qualifier verification are included below.

2.1 SDG JC7512

Data validation and review was conducted for soil samples within SDG JC7512 in accordance with the EPA National Functional Guidelines for Superfund Organic Methods Review, January 2017, EPA National Functional Guidelines for Superfund Inorganic Methods Review, January 2017, and NYDEC Analytical Services Protocol. The SDG included six soil samples to be analyzed for VOCs by EPA 8260C, SVOCs by EPA 8270D, chlorinated herbicides by EPA 8151, chlorinated pesticides by EPA 8081B, polychlorinated biphenyls by EPA 8082A, metals by EPA 6010C, mercury by EPA 7471B and percent moisture by SM2540 G-97. Samples specific to JC7512 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|----------|-----------|--------|-------------|-------------|
| JC7512-1 | SB-1-1 | Soil | 10/29/2015 | 13:00 |
| JC7512-2 | SB-1-5 | Soil | 10/29/2015 | 13:10 |
| JC7512-3 | SB-2-1 | Soil | 10/29/2015 | 13:20 |
| JC7512-4 | SB-2-5 | Soil | 10/29/2015 | 13:55 |
| JC7512-5 | SB-3-1 | Soil | 10/29/2015 | 14:30 |
| JC7512-6 | SB-3-5 | Soil | 10/29/2015 | 14:45 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: the cooler temperature upon receipt was 2.5°C, custody seals were intact and containers were accurately labeled and reflected the chain of custody. Soil samples were preserved onsite with DI water and methanol for VOC analysis. Samples preserved with DI water were frozen by the analytical laboratory upon receipt.

2.1.1 VOCs

Holding Times

Samples were received below 6°C and within method holding times by the analytical laboratory. Soil samples preserved with DI water were frozen upon receipt by the analytical laboratory. Analysis for VOCs was performed between five and eight days from sample collection. Analysis times were confirmed with raw data.

Instrument Performance Check

Bromofluorobenzene (BFB) was evaluated for ion abundance requirements and frequency to ensure adequate instrument performance. Four sequences and subsequently four BFB standards were analyzed for samples included in SDG JC7512.

The initial sequence, with an injection date of 9/17/15, includes the soil matrix calibration and initial calibration check. BFB was analyzed at the beginning of the sequence and within ion abundance criteria. The final sample for the sequence was analyzed six hours past the BFB injection. The BFB ion abundances were reviewed from raw data presented in the SDG. Ion abundances were averaged over three scans and background-subtracted from ten scans away from the BFB peak.

Three more BFB injections were performed. One BFB injection was completed on 11/2/15 for JC7512-1, JC7512-2 and quality control samples. The ion abundances were averaged over three scans and background subtraction was performed ten scans away from the BFB peak. The second BFB injection was completed on 11/3/15 for JC7512-4, JC7512-5, JC7512-6 and quality control samples. Ion abundances were averaged over three scans and were background subtracted from ten scans away from the BFB peak. The final BFB injection was performed on 11/6/15 for JC7512-3 and quality control samples. The ion abundances were averaged over three scans and were background subtracted from ten scans away from the BFB peak. The final BFB injection was performed on 11/6/15 for JC7512-3 and quality control samples. The ion abundances were averaged over three scans and were background subtracted from ten scans away. All BFB injections were within twelve hours of sample injections and were within ion abundance criteria.

Initial Calibration

Initial calibration included at least five calibration points, were performed prior to ICVs, samples and blanks, and were within twelve hours of a BFB check standard. Relative Response Factors (RRF) and Percent Relative Standard Deviations (%RSD) were calculated for each analyte and surrogate. The mean RRF and %RSD for target analytes were within control limits as stated in Table 18 of the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017.

Initial Calibration Verification

The RRF for target analytes and surrogate analytes were above minimums presented in the CLP Guidance. The percent difference between the mean RRF of the initial calibration and the initial calibration verification were within control limits stated by the CLP Guidance.

Continuing Calibration Verification

Continuing calibration verifications were evaluated to determine instrument performance and calibration validity during each analytical run. The RRF calculated for each analyte and the %D from the initial calibration were within control limits except for dichlorodifluoromethane from the CCV analyzed 11/2/15 at 10:37 pm and 11/6/15 at 10:32 am. Samples within analytical sequences affected by this %RSD failure are JC7512-1 (SB-1-1), JC7512-2 (SB-1-5) and JC7512-3 (SB-2-1). Dichlorodifluoromethane is estimated for these samples and is UJ qualified within the data table.

Blanks

Three method blanks, one per analytical batch, were analyzed to assess possible contamination from the laboratory or the field. There were no detections in any blanks associated with this SDG. Chromatograms were checked to ensure that all measurable peaks were identified.

Surrogates

The percent recovery of surrogate compounds were evaluated to determine method and instrument performance. Surrogate recoveries for soil samples and quality control samples were within laboratory control limits. Raw data was checked for calculation and transcription errors and none were found.

Laboratory Control Sample

One laboratory control sample was analyzed per analytical batch included in the SDG. The percent recoveries for target analytes were within laboratory control limits. Raw data was compared to reported results and no

discrepancies were found.

Matrix Spike/Matrix Spike Duplicate

One matrix spike was analyzed for samples JC7512-4 (SB-2-5), JC7512-5 (SB-3-1) and JC7512-6 (SB-3-5). The recovery of MS analytes was within laboratory and CLP control limits. Two analytical batches included MS/MSD analysis. Both had multiple target analytes outside of percent recovery and RPD control limits. Samples used as the source for these MS/MSDs were not part of SDG JC7512 and no data was qualified in the data table.

Laboratory Duplicate

A laboratory duplicate was performed on sample JC7512-5 (SB-3-1). RPD values were within laboratory control limits.

Internal Standard

Required internal standards were added to samples, blank samples and quality control samples at the same specified concentration to monitor system performance and quantify target analytes. Raw data was checked against acceptance criteria for retention times and area response. Retention times did not differ more than 10 seconds from the check standard (beginning calibration check verification). Area responses were within 50-200% of the check standard. Raw data was compared to reported summary results and no discrepancies were found.

Target Analyte Identification

Target analyte detections are confirmed by ion fragments and retention times. Acetone had secondary and tertiary ion fragments outside of the acceptable abundance range as defined by the initial calibration. These results were accepted as positively identified compounds based on RT and analyte mass spectra compared to the reference spectra. Trichloroethene ion abundances were also outside of acceptance criteria but this is most likely due to the analyte concentration being below the reporting limit. Both retention times and comparison to the reference spectrum suggest that TCE is present within samples.

Target Analyte Quantitation and Reported Contract Required Quantitation Limits

Results for positively identified analytes and method reporting limits were calculated correctly by the laboratory and adjusted based on initial sample volumes and percent solids. The percent solids in the soil samples were greater than 30%. Internal standards based on the ICAL were used for quantitation of analytes in samples and quality control samples. Primary ion fragments used to identify analytes and mean RRF were the same as the ICAL.

2.1.2 SVOCs

Preservation and Holding Times

Samples were received by the laboratory below 6°C, were extracted five days from sample collection and analyzed the same day as extraction. Extraction dates and analysis dates from raw data were confirmed against reported dates.

GC/MS Instrument Performance Check

Decafluorotriphenylphosphine (DFTPP) mass spectra and mass listing was evaluated for ion abundance requirements and analysis frequency to ensure adequate instrument performance. Raw data was evaluated to ensure instrument tuning was within CLP requirements. DFTPP was analyzed at the beginning of each sequence and within 12 hours of the last sample injection. Raw data abundances and calculations were checked against tune reports. No discrepancies were found.

Initial Calibration

Initial calibration was performed before ICV, samples and required blanks and within 12 hours of the DFTPP tune

check. A minimum of five standards that ranged between 1 ppm and 100 ppm were used to calibrate the instrumentation. The initial calibration summary report was checked for correct calculation of RRF, mean RRF and %RSD. The initial calibration summary report was generated from acquisition software. RRF and %RSD of target analytes were within CLP guidance except for 2,4-dinitrophenol which was calibrated using a quadratic regression with a correlation coefficient of 0.999.

Initial Calibration Verification

Initial calibration verification was performed after calibration and before any samples, quality control samples, blank analysis and continuing calibration verification. The ICV standards were prepared at 50 ppm, which is the middle of the calibration curve. The RRF for target analytes were above minimum RRFs stated in CLP guidance. The %RSD for 2-nitroaniline and carbazole were outside of control limits and are UJ-flagged in the data table.

Continuing Calibration Verification

CCV was analyzed at the mid-point of the calibration curve. RRF of analytes were compared to mean RRF from the initial calibration, all were within CLP control limits. The %RSD between the ICAL mean RRF and the CCV target analyte RRF were outside of CLP control limits for nitrobenzene-d5 and hexachlorobutadiene. Results for the DMC nitrobenzene-d5 should be considered estimated. Hexachlorobutadiene is UJ-flagged in the data table as an estimated value.

Blanks

One method blank was extracted with the associated soil samples in SDG JC7512. There were no detections of target analytes within the method blank. The method blank chromatogram was also reviewed to ensure all peaks were identified.

Surrogates

The percent recovery of surrogate compounds was within CLP and laboratory control limits. Each surrogate was accurately identified, as confirmed after raw data review of retention times and primary ion.

Laboratory Control Sample

One laboratory control sample was extracted and analyzed with samples of SDG JC7512. Percent recoveries were within control limits.

Matrix Spike/Matrix Spike Duplicate

One matrix spike and matrix spike duplicate pair were extracted and analyzed with samples from SDG JC7512. The source sample used for the MS/MSD was JC7512-1 and percent recoveries and RPD values of target analytes were within control limits.

Internal Standard

Internal standards were added to all samples, quality control samples and blanks. Area and retention times for internal standards were within limits set by the check standard. The check standard used was the initial CCV on 11/3/15. Raw data was compared to summary data pages and no discrepancies were found.

Target Analyte Identification

The RRT of bis(2-ethylhexyl)phthalate was greater than the 0.06 RT units control limit from sample JC7512-3 (SB-2-1). Due to the mass spectra and ion abundance criteria within control limits, identification of the analyte is accepted.

Target Analyte Quantitation and Reported Contract Require Quantitation Limit

Sample results and reporting limits were correctly calculated based on sample preparation initial and final volumes and the percent solids within the soil samples. Correct internal standards were used to quantitate target analytes. The same mean RRF and quantitation ions were used consistently. Soil samples had greater than 30% total solids.

2.1.3 Pesticides

Preservation and Holding Times

Samples were extracted within five days of samples collection and were analyzed two days after extraction which meets method holding requirements. Samples were not chemically preserved but were held by the laboratory below 6°C.

GC/ECD Performance Check

The breakdown of DDT and Endrin was below 20%. Breakdown checks were analyzed at the beginning of every sequence and within 12 hours of the final sample injected into the instrument.

Initial Calibration

Two calibrations were performed to quantitate soil and quality control samples. Target analytes were calibrated at eight concentration levels. A single point toxaphene and chlordane standard were included. Raw data was checked to verify concentration levels were accurately reported. Mean retention times were calculated from three check standards. Response factors were calculated for target analytes and %RSDs were below 20% except for 4,4'-DDT on column 2 of calibration ICC840. Calibration was accepted based on the quadratic calibration for 4,4'-DDT with a correlation coefficient of 0.9996.

Initial and Continuing Calibration

The %D between the average RF from the ICAL and the initial and continuing calibration checks were within the $\pm 25\%$ range. Retention times were within windows.

Blanks

A method blank was extracted in a batch of nine samples that included samples from SDG JC7512. Instrument blanks were analyzed after initial and continuing calibration verification. An erroneous peak was found in the method blank, but retention times did not confirm it with any target analytes.

Surrogate

TCX and DCB were added to all SDG samples and quality control samples prior to extraction. Retention times for the surrogate compounds were within RT windows. Surrogate recoveries were within laboratory control limits.

Laboratory Control Sample

One laboratory control sample was prepared with samples from SDG JC7512 on 11/3/15. The recovery of target analytes were within CLP guidance limits.

Matrix Spike/Matrix Spike Duplicate

A sample not included in SDG JC7512 was used as the source for the MS/MSD. The RPD value was exceeded based on CLP guidance for dieldrin. Since the source sample was not part of JC7512, no sample data was qualified.

Target Analyte Identification

4,4'-DDD was detected in JC7512-1 (SB-1-1) but the RPD between the concentrations on the two columns exceeded 25%. This analyte has been updated to its DL and qualified with a U in the data table.

2.1.4 PCBs

Preservation and Holding Times

Samples were extracted within five days of samples collection and were analyzed three to four days after extraction which is within method holding times. Samples were not chemically preserved but were stored by the laboratory below 6°C.

Initial Calibration

A five-point calibration curve was generated for Aroclor 1016 and 1260. A one-point calibration curve was generated for Aroclors 1221, 1232, 1242, 1248, 1254, 1262 and 1268. Mean retention times were calculated from the 1000 ppb ICAL standard and two subsequent continuing calibration verifications. The RF was calculated for target peaks and the %RSD between the calibration levels was less than 20%.

Continuing Calibration Verification

The %D between the ICAL mean RF and the ICV/CCV were within ±35%. All Aroclor peaks were within RT windows set by the ICAL. Instrument blanks bracketed CCVs and were analyzed less than 12 hours apart.

Blanks

A method blank was included in the extraction batch. No Aroclor patterns were found.

Surrogate

A surrogate mixture of TCX and DCM was added to all samples and quality control samples prior to extraction. The recovery of the surrogate compounds TCX and DCM were between 30-150% and retention time shifts were within windows.

Matrix Spike/Matrix Spike Duplicate

One matrix spike and matrix spike duplicate pair were extracted in the batch that included samples in SDG JC7512. The source for the matrix spike and matrix spike duplicate samples was not part of SDG JC7512. Spike recovery and RPD values between MS and MSD were within control limits and were accurately calculated.

Laboratory Control Sample

A LCS was extracted and analyzed with samples from SDG JC7512. Raw data was verified against report lab values by recalculation. The percent recovery of spiked Aroclor 1016 and 1260 were within control limits of 50-150%.

2.1.5 Herbicides

Preservation and Holding Times

Samples were extracted within five days of samples collection and were analyzed seven days after extraction.

Initial Calibration

A six-point calibration curve was generated for 2,4-D, 2,4,5-TP and 2,4,5-T. The RF was calculated for target peaks and the %RSD between the calibration levels was within ±20%. Mean retention times and RT windows were calculated from ICV and CCV after the ICAL. Analyte identification is based off these RT windows.

Continuing Calibration Verification

The %D between the ICAL mean RF and the ICV/CCV were within ±35% for target analytes. All target analyte peaks were within RT windows.

Blanks

A method blank was included in the extraction batch. No target analytes were found. Raw data was evaluated for any erroneous analyte peaks. Peaks found did not have retention times that aligned with target analytes.

Surrogate

2,4-DCAA was used as the surrogate for all sample and quality control samples analyzed for herbicides. Low recovery of surrogate was found in all soil samples and MS/MSDs extracted, but laboratory control limits were not exceeded. Target analyte recovery might be compromised due to matrix. Retention times were verified and compared to the check standard.

Matrix Spike/Matrix Spike Duplicate

One matrix spike and matrix spike duplicate pair were extracted in the batch that included samples in SDG JC7512. The source for the matrix spike and matrix spike duplicate samples was JC7512-1. Spike recovery and RPD values between MS and MSD were within control limits and were accurately calculated. The retention time of analytes were within windows.

Laboratory Control Sample

A LCS was extracted and analyzed with samples from SDG JC7512. The percent recovery of spiked target analytes was within control limits.

2.1.6 Metals

Preservation and Holding Time

Samples were prepared and analyzed within one week of arrival at the laboratory. Preparation and analysis dates that are reported were confirmed with raw data.

Calibration

Target analytes were calibrated using a 2-point calibration curve. The initial and continuing calibration verification standards were within 10% of their true values.

Blanks

A CCB was analyzed after every CCV, which was analyzed every 10 samples. The value of CCB did not exceed laboratory reporting limits. The method blank did not exceed laboratory reporting limits.

Interference Check Standard

The ICSAB and ICSA check standards were within 20% of the true value and ND values were no greater than \pm RL.

Laboratory Control Sample

Target analyte recoveries were within ±30% of true values.

Matrix Spike/Matrix Spike Duplicate

The source for the MS/MSD is not included in SDG JC7512; a serial dilution was also performed on this sample. The recovery of aluminum was above the upper control limit. No data was qualified since the source sample was not included in SDG JC7512.

2.1.7 Mercury

Preservation and Holding Times

Samples were prepared six days from collection and analyzed within 24 hours. Raw data was reviewed to verify sample preparation and analysis dates.

Calibration

The calibration curve consisted of 5 points with a correlation coefficient of 0.999. CCV4 has a recovery of mercury at 116%. Prep batch quality control may be biased high. Sample data was not affected.

Blanks

The CCB and method blanks were below the reporting limit.

Spiked Samples

The laboratory control sample was within $\pm 20\%$ of the true value of the mercury spike. The MS/MSD were not prepared with a sample from SDG JC7512, but all recoveries were within $\pm 20\%$ and within the RPD limit of 20%.

2.1.8 Percent Moisture

Percent moisture was performed and calculated for soil samples included in SDG JC7512. Samples results were verified by recalculation of raw data, where; $Percent\ Moisture = \frac{Wet\ Weight-Dry\ weight}{Wet\ Weight-Tare\ Weight}$ x100%. Results reported by the laboratory are calculated on a dry weight basis.

2.2 SDG JC7519

Data validation and review was conducted for air samples within SDG JC7519 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017, EPA method TO-15 and NYDEC Analytical Services Protocol. The SDG included three air samples to be analyzed for VOCs by TO-15. Samples specific to JC7519 are included below.

| SDG ID Sample ID | | Matrix | Sample Date | Sample Time |
|------------------|-------|--------|-------------|-------------|
| JC7519-1 | SV-09 | Air | 10/29/2015 | 8:55 |
| JC7519-2 | SV-10 | Air | 10/29/2015 | 8:45 |
| JC7519-3 | SV-11 | Air | 10/29/2015 | 8:30 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: custody seals were intact and containers were accurately labeled and reflected the chain of custody. The air samples were collected in 6-liter canisters and no damage was present when received by the laboratory.

Preservation and Holding Time

The air samples were collected in cleaned metal canisters. They arrived at the analytical laboratory two days after sampling and were analyzed between three and five days from receipt. Raw data was verified to ensure analysis date reported is correct.

GC/MS Instrument Performance Check

BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence, per TO-15. Ion abundances were averaged over three scans and background was subtracted from less than 20 scans away from the BFB peak.

Initial Calibration

RRF and %RSD for target analytes were within control limits for all calibrations associated with SDG JC7519. A minimum of 5 calibration points were used for the calibration curves.

Initial Calibration Verification

An ICV standard was analyzed after each ICAL and before any samples or quality control samples. The ICV following calibration on 10/5/15 had a RRF value for 1,2,4-trichlorobenze below the minimum value of 0.300. No samples or quality control samples included in analysis for SDG JC7519 are affected by this since the calibration was for Summa Cleaning Certification samples.

Continuing Calibration Verification

Target analytes were within %D limits for opening CCVs and above minimum RRFs as stated by CLP Guidance.

Blanks

Method blanks were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. No analytes other than surrogates and internal standards were detected in method blanks. Raw data was verified that no other peaks were present in the chromatograms.

Surrogates

The surrogate was within laboratory control limits. Raw data for samples was checked to ensure that when dilutions were run, both recoveries were within control limits.

Laboratory Control Samples

LCS/LCSDs were performed with each sequence where analysis of environmental samples or SCC samples were performed. Recoveries and RPD values of target analytes were within control limits.

Laboratory Duplicate

A laboratory duplicate was performed on a sample not included in SDG JC7519. All RPD values were within control limits.

Summa Cleaning Certification

Summa's were cleaned and no residual VOCs were found.

Internal Standards

The retention time of the internal standards was compared against daily continuing calibration verifications. The retention times did not differ more than 0.33 minutes and area counts were within 40% of the check standard for samples associated with JC7519.

Target Analyte Identification

Samples associated with JC7519 had complicated matrices with a mixture of different VOCs. Raw data was verified to ensure that analyte identification was accurate by comparing ion abundances and retention times to check standards. Some target analytes did not meet the abundance criteria of ±20% due to matrix interference or low

analyte concentration. These analytes were reported by the laboratory based on RT and analyte mass spectra.

2.3 SDG JC7667

Data validation and review was conducted for air samples within SDG JC7667 in accordance with the EPA National Functional Guidelines for Superfund Organic Methods Review, January 2017, EPA Method TO-15 and NYDEC Analytical Services Protocol. The SDG included three air samples that were analyzed for VOCs by TO-15. Samples specific to JC7667 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|----------|-----------|--------|----------------|-------------|
| JC7667-1 | SV-8 | Air | 10/30/2015 | 6:55 |
| JC7667-2 | SV-7 | Air | 10/30/2015 | 7:05 |
| JC7667-3 | SV-6 | Air | 10/30/2015 | 11:10 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey on November 3, 2015. As indicated by the analytical laboratory sample receipt summary: custody seals were intact and containers were accurately labeled and reflected the chain of custody. The air samples were collected in 6-liter canisters and no damage was present when received by the laboratory.

Preservation and Holding Time

The air samples were collected in cleaned metal canisters and were analyzed within method holding limits. They arrived at the analytical laboratory four days after sampling and were analyzed on November 5, 2015 and November 6, 2015. Raw data was verified to ensure analysis date reported is correct.

GC/MS Instrument Performance Check

BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence and 24 hours within the last samples injected, per method TO-15. Ion abundances were averaged over three scans and were background subtracted from less than 20 scans away.

Initial Calibration

RRF and %RSD were within control limits for all calibrations associated with SDG JC7667. A minimum of 5 calibration points were used for the calibration curves.

Initial Calibration Verification

An ICV standard was analyzed after each ICAL. Minimum RRF was met for target analytes and %D between the mean ICAL RRF and the ICV RRF was within control limits.

Continuing Calibration Verification

Target analytes were within %D limits for opening CCVs.

Blanks

Method blanks were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. No analytes other than surrogates and internal standards were detected in method blanks. Raw data was verified that no other peaks were present in the chromatograms.

Surrogates

The recovery of the surrogate compound was within control limits for samples and associated quality control samples within the SDG.

Laboratory Control Samples

LCS/LCSDs were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. Recoveries and RPD values of target analytes were within control limits.

Laboratory Duplicate

The source for the laboratory duplicate analyzed was not part of SDG JC7667. All RPD values were <20%.

Summa Cleaning Certification

Summa's were cleaned and no residual VOCs were found.

Internal Standards

The retention time of the internal standards was compared against daily opening continuing calibration verifications. The retention times did not differ more than 0.33 minutes and area counts were within 40% of the check standard.

Target Analyte Identification

Raw data was verified to ensure that analyte identification was accurate by comparing ion abundances, analyte mass spectra and retention times to check standards. Some target analytes did not meet the abundance criteria of $\pm 20\%$ due to matrix interferences but RT and mass spectra suggest correct analyte identification.

2.4 SDG JC12861

Data validation and review was conducted for soil and blank water samples within SDG JC12861 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 and NYDEC Analytical Services Protocol. The SDG included eight soil samples, one trip blank and one field blank to be analyzed for VOCs by EPA 8260C and percent moisture by SM2540. Samples specific to JC12861 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|------------|----------------|------------------|-------------|-------------|
| JC12861-1 | MW-11-20' | Soil | 1/18/2016 | 10:20 |
| JC12861-2 | MW-11-45' | Soil | 1/18/2016 | 10:55 |
| JC12861-3 | MW-11-50' | Soil | 1/18/2016 | 11:10 |
| JC12861-4 | MW-1I | Soil | 1/18/2016 | 12:00 |
| JC12861-5 | Trip | Trip Blank (aq) | 1/18/2016 | |
| JC12861-6 | Field | Field Blank (aq) | 1/19/2016 | 11:42 |
| JC12861-7 | MW-45-23.5' | Soil | 1/19/2016 | 10:20 |
| JC12861-8 | MW-45-30' | Soil | 1/19/2016 | 10:30 |
| JC12861-9 | MW-45-30' | Soil | 1/19/2016 | 10:50 |
| JC12861-10 | MW-45-MOISTURE | Soil | 1/19/2016 | 13:20 |

Samples were received within twenty-four hours of sampling by the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: the cooler temperature upon receipt was 1.3°C, custody seals were intact, containers were accurately labeled and reflected the chain of custody and there was no headspace in VOA (volatile organic analysis) water containers. Soil samples were collected using field core sampling with enclosed storage containers (EnCore sampler) in compliance with preparation method EPA 5035 and were prepared by the analytical laboratory within twenty-four hours of sample receipt. The QC blank samples were preserved with hydrochloric acid (HCl) and pH checked by the laboratory.

2.4.1 VOCs

Holding Times

Samples were received below 6°C and within method holding limits by the analytical laboratory. Soil samples were prepared by the laboratory within twenty-four hours of receipt. QC blank samples were preserved with HCl and tested below a pH of 2 by the laboratory. There was no headspace in VOA containers. Analysis for VOCs was performed between three and six days from sample collection. Holding times for soil and water samples analyzed for VOCs were met.

Instrument Performance Check

Bromofluorobenzene (BFB) was evaluated for mass spectra requirements and frequency to ensure adequate instrument performance. Four sequences and subsequently four BFB standards were analyzed for samples included in SDG JC12861.

The initial sequence with injection dates of 1/12/16 includes the water matrix calibration and calibration check. BFB was analyzed at the beginning of the sequence and within ion abundance criteria. The final sample for the sequence was analyzed six hours and forty minutes past the BFB injection. The BFB ion abundances were reviewed from raw data presented in the SDG, recalculated to ensure data validity and presented below. Ion abundances were averaged over three scans and background-subtracted from eleven scans away.

BFB Injection: 1/12/16 14:11

| 36.00 717 37.10 3783 38.10 3372 39.05 1194 39.05 1194 39.05 1194 39.05 1194 39.05 1194 39.05 1194 39.05 1194 39.05 1194 39.05 62 44.00 332 45.10 569 47.05 832 48.00 427 49.05 3385 50.10 14270 51.05 4365 52.05 165 55.10 66 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3353 62.05 3025 63.00 2476 64.10 198 67.15 214 | m/z | abund. |
|---|-------|--------|
| 38.10 3372 39.05 1194 39.90 62 44.00 332 45.10 569 47.05 832 48.00 427 49.05 3385 50.10 14270 51.05 4365 52.05 165 55.10 66 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3325 62.05 3025 63.00 2476 64.10 198 | 36.00 | 717 |
| 39.05 1194 39.90 62 44.00 332 45.10 569 47.05 832 48.00 427 49.05 3385 50.10 14270 51.05 4365 52.05 165 55.10 66 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3325 62.05 3025 63.00 2476 64.10 198 | 37.10 | 3783 |
| 39.90 62 44.00 332 45.10 569 47.05 832 48.00 427 49.05 3385 50.10 14270 51.05 4365 52.05 165 55.10 66 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3325 62.05 3025 63.00 2476 64.10 198 | 38.10 | 3372 |
| 44.0033245.1056947.0583248.0042749.05338550.101427051.05436552.0516555.106656.05109457.05200858.107060.0565361.05335362.05302563.00247664.10198 | 39.05 | 1194 |
| 45.1056947.0583248.0042749.05338550.101427051.05436552.0516555.106656.05109457.05200858.107060.0565361.05335362.05302563.00247664.10198 | 39.90 | 62 |
| 47.0583248.0042749.05338550.101427051.05436552.0516555.106656.05109457.05200858.107060.0565361.05335362.05302563.00247664.10198 | 44.00 | 332 |
| 48.0042749.05338550.101427051.05436552.0516555.106656.05109457.05200858.107060.0565361.05335362.05302563.00247664.10198 | 45.10 | 569 |
| 49.05338550.101427051.05436552.0516555.106656.05109457.05200858.107060.0565361.05335362.05302563.00247664.10198 | 47.05 | 832 |
| 50.101427051.05436552.0516555.106656.05109457.05200858.107060.0565361.05335362.05302563.00247664.10198 | 48.00 | 427 |
| 51.05 4365 52.05 165 55.10 66 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3353 62.05 3025 63.00 2476 64.10 198 | 49.05 | 3385 |
| 52.05 165 55.10 66 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3353 62.05 3025 63.00 2476 64.10 198 | 50.10 | 14270 |
| 55.10 66 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3353 62.05 3025 63.00 2476 64.10 198 | 51.05 | 4365 |
| 56.05 1094 57.05 2008 58.10 70 60.05 653 61.05 3353 62.05 3025 63.00 2476 64.10 198 | 52.05 | 165 |
| 57.05200858.107060.0565361.05335362.05302563.00247664.10198 | 55.10 | 66 |
| 58.10 70 60.05 653 61.05 3353 62.05 3025 63.00 2476 64.10 198 | 56.05 | 1094 |
| 60.0565361.05335362.05302563.00247664.10198 | 57.05 | 2008 |
| 61.05335362.05302563.00247664.10198 | 58.10 | 70 |
| 62.05302563.00247664.10198 | 60.05 | 653 |
| 63.00247664.10198 | 61.05 | 3353 |
| 64.10 198 | 62.05 | 3025 |
| | 63.00 | 2476 |
| 67.15 214 | 64.10 | 198 |
| | 67.15 | 214 |

| m/z | abund. |
|--------|--------|
| 68.00 | 7386 |
| 69.05 | 7372 |
| 70.05 | 660 |
| 72.00 | 459 |
| 73.00 | 3353 |
| 74.00 | 12378 |
| 75.05 | 35056 |
| 76.00 | 3165 |
| 77.05 | 470 |
| 78.10 | 102 |
| 78.90 | 1894 |
| 80.00 | 605 |
| 80.90 | 2004 |
| 81.95 | 456 |
| 87.00 | 2798 |
| 87.95 | 2686 |
| 90.95 | 279 |
| 92.00 | 1925 |
| 93.00 | 3205 |
| 94.10 | 8775 |
| 95.00 | 70125 |
| 96.00 | 4663 |
| 103.80 | 107 |

| m/z | abund. |
|--------|--------|
| 104.00 | 217 |
| 104.90 | 67 |
| 105.95 | 331 |
| 115.95 | 314 |
| 116.95 | 479 |
| 117.95 | 286 |
| 118.85 | 414 |
| 127.95 | 261 |
| 129.00 | 72 |
| 129.95 | 334 |
| 130.90 | 76 |
| 140.95 | 857 |
| 142.95 | 896 |
| 148.00 | 66 |
| 154.90 | 147 |
| 158.90 | 60 |
| 171.70 | 72 |
| 171.90 | 68 |
| 173.95 | 71298 |
| 175.00 | 6036 |
| 175.95 | 69288 |
| 176.95 | 4528 |
| | |

| Mass | Criteria | % Rel. Abundance |
|------|-----------------------------------|------------------|
| 50 | 15-40% of mass 95 | 20.35% |
| 75 | 30-80% of mass 95 | 49.99% |
| 95 | Base peak, 100% of rel. abundance | 100.00% |
| 96 | 5-9% of mass 95 | 6.65% |
| 173 | <2% of mass 174 | 0.00% |
| 174 | 50-120% of mass 95 | 101.67% |
| 175 | 5-9% of mass 174 | 8.47% |
| 176 | 95-101% of mass 174 | 97.18% |
| 177 | 5-9% of mass 176 | 6.54% |

A second injection of BFB was performed on 1/21/16 which encompassed quality control samples for a water matrix analytical batch along with the Trip Blank (JC12861-5) and Field Blank (JC12861-6) samples. The BFB sample was analyzed at the beginning of the sequence and was within ion abundance criteria. Abundances were averaged over three scans and background was subtracted from thirteen scans away. The BFB standard was injected ten hours and

forty-six minutes before the final sample in the sequence. Raw data and calculated abundance criteria are present below.

BFB Injection: 1/27/16 21:21

| m/z | abund. |
|-------|--------|
| 36.05 | 669 |
| 37.10 | 3355 |
| 38.10 | 2994 |
| 39.10 | 1216 |
| 44.00 | 239 |
| 45.05 | 634 |
| 47.05 | 594 |
| 48.05 | 465 |
| 49.05 | 2859 |
| 50.10 | 12814 |
| 51.10 | 3686 |
| 52.05 | 154 |
| 55.00 | 138 |
| 56.00 | 1013 |
| 57.05 | 1942 |
| 60.05 | 729 |
| 61.05 | 3200 |
| 62.05 | 3081 |
| 63.00 | 2193 |
| 64.10 | 136 |
| 68.00 | 6904 |
| 69.05 | 6755 |
| 70.10 | 484 |

| m/z | abund. |
|--------|--------|
| 72.05 | 400 |
| 73.05 | 2907 |
| 74.05 | 11263 |
| 75.00 | 32347 |
| 76.05 | 2770 |
| 76.95 | 346 |
| 77.95 | 184 |
| 78.90 | 1959 |
| 80.00 | 576 |
| 81.00 | 1940 |
| 81.90 | 374 |
| 87.00 | 2122 |
| 88.00 | 2110 |
| 90.80 | 61 |
| 90.95 | 242 |
| 92.00 | 1844 |
| 93.00 | 2839 |
| 94.05 | 7639 |
| 95.00 | 62933 |
| 96.05 | 4343 |
| 104.00 | 407 |
| 104.90 | 68 |
| 105.95 | 23 |

| m/z | abund. |
|--------|--------|
| 115.95 | 233 |
| 117.00 | 461 |
| 117.95 | 283 |
| 118.95 | 412 |
| 127.90 | 270 |
| 128.90 | 84 |
| 129.95 | 299 |
| 134.90 | 68 |
| 136.90 | 72 |
| 140.95 | 876 |
| 141.90 | 65 |
| 142.95 | 909 |
| 147.85 | 258 |
| 154.90 | 189 |
| 156.80 | 61 |
| 171.30 | 68 |
| 171.90 | 75 |
| 172.10 | 98 |
| 174.00 | 63291 |
| 175.00 | 5224 |
| 175.95 | 60696 |
| 176.95 | 4007 |
| 177.95 | 219 |
| • | • |

| Mass | Criteria | % Rel. Abundance |
|------|-----------------------------------|------------------|
| 50 | 15-40% of mass 95 | 20.36% |
| 75 | 30-80% of mass 95 | 51.40% |
| 95 | Base peak, 100% of rel. abundance | 100.00% |
| 96 | 5-9% of mass 95 | 6.90% |
| 173 | <2% of mass 174 | 0.00% |
| 174 | 50-120% of mass 95 | 100.57% |
| 175 | 5-9% of mass 174 | 8.25% |
| 176 | 95-101% of mass 174 | 95.90% |
| 177 | 5-9% of mass 176 | 6.60% |

For soil analysis, BFB was analyzed at the beginning of the calibration on 1/5/16. The BFB standard was injected six hours and nine minutes before the final sample in the sequence. The ion abundances were within guidance criteria, averaged over three scans and background subtracted from eleven scans away. Raw data and calculated abundance criteria are presented below.

BFB Injection: 1/5/16 9:36

| m/z | abund. |
|-------|--------|
| 36.00 | 959 |
| 37.00 | 5017 |
| 38.00 | 4536 |
| 39.00 | 1665 |
| 41.00 | 121 |
| 42.00 | 127 |
| 43.00 | 113 |
| 44.00 | 359 |
| 45.00 | 970 |
| 46.00 | 58 |
| 47.00 | 1348 |
| 48.00 | 663 |
| 49.05 | 4130 |
| 50.00 | 19080 |
| 51.00 | 5875 |
| 52.05 | 327 |
| 55.00 | 157 |
| 56.00 | 1463 |
| 57.00 | 2620 |
| 58.10 | 65 |
| 60.00 | 922 |
| 61.00 | 4720 |
| 62.00 | 4248 |
| 63.00 | 3298 |
| 64.05 | 307 |
| 67.05 | 416 |
| 68.00 | 10176 |
| | |

| m/z | abund. |
|--------|--------|
| 69.00 | 10013 |
| 69.95 | 781 |
| 72.00 | 438 |
| 73.00 | 4225 |
| 74.00 | 16637 |
| 75.00 | 48354 |
| 76.00 | 4177 |
| 76.95 | 648 |
| 77.90 | 345 |
| 78.10 | 180 |
| 78.90 | 2763 |
| 79.95 | 791 |
| 80.90 | 2962 |
| 81.90 | 679 |
| 87.00 | 4620 |
| 87.95 | 4232 |
| 90.90 | 485 |
| 92.00 | 2601 |
| 93.00 | 4052 |
| 94.00 | 11273 |
| 95.00 | 100213 |
| 96.00 | 6769 |
| 97.00 | 165 |
| 103.95 | 496 |
| 104.80 | 87 |
| 105.85 | 424 |
| 111.80 | 59 |

| | 1 |
|--------|--------|
| m/z | abund. |
| 115.95 | 345 |
| 116.95 | 645 |
| 117.90 | 387 |
| 118.95 | 558 |
| 127.85 | 400 |
| 128.85 | 135 |
| 129.95 | 342 |
| 130.95 | 194 |
| 134.90 | 291 |
| 136.85 | 208 |
| 139.70 | 60 |
| 140.90 | 920 |
| 141.80 | 58 |
| 142.95 | 947 |
| 146.00 | 66 |
| 147.00 | 45 |
| 147.60 | 64 |
| 148.00 | 77 |
| 154.80 | 144 |
| 156.80 | 65 |
| 157.00 | 90 |
| 173.90 | 78530 |
| 174.95 | 5936 |
| 175.90 | 76752 |
| 176.90 | 5421 |
| 177.90 | 68 |
| 253.05 | 8 |

| Mass | Criteria | % Rel. Abundance |
|------|-----------------------------------|------------------|
| 50 | 15-40% of mass 95 | 19.04% |
| 75 | 30-80% of mass 95 | 48.25% |
| 95 | Base peak, 100% of rel. abundance | 100.00% |
| 96 | 5-9% of mass 95 | 6.75% |
| 173 | <2% of mass 174 | 0.00% |
| 174 | 50-120% of mass 95 | 78.36% |
| 175 | 5-9% of mass 174 | 7.56% |
| 176 | 95-101% of mass 174 | 97.74% |
| 177 | 5-9% of mass 176 | 7.06% |

A final BFB standard was analyzed for soil samples and guidance criteria was met. The BFB standard was injected ten hours and thirty minutes before the final sample in the sequence. The ion abundances were averaged over three scans and background subtracted from eleven scans away. Raw data and calculated abundance criteria are presented below.

BFB Injection: 1/24/16 10:41

| m/z | abund. |
|-------|--------|
| 36.10 | 1412 |
| 37.10 | 7945 |
| 38.05 | 6897 |
| 39.10 | 2640 |
| 40.00 | 183 |
| 43.05 | 130 |
| 44.00 | 783 |
| 45.05 | 1420 |
| 47.05 | 2025 |
| 48.00 | 924 |
| 49.05 | 6334 |
| 50.10 | 30901 |
| 51.10 | 9233 |
| 52.05 | 453 |
| 55.05 | 372 |
| 56.00 | 2164 |
| 57.00 | 4198 |
| 57.95 | 205 |
| 60.05 | 1388 |
| 61.00 | 6710 |
| 62.05 | 6703 |
| 63.00 | 4908 |

| m/z | abund. | | | | |
|--------|--------|--|--|--|--|
| 73.00 | 6178 | | | | |
| 74.00 | 23746 | | | | |
| 75.00 | 71661 | | | | |
| 76.05 | 6275 | | | | |
| 76.95 | 1000 | | | | |
| 78.05 | 805 | | | | |
| 78.90 | 4082 | | | | |
| 79.95 | 1335 | | | | |
| 80.90 | 4039 | | | | |
| 81.95 | 936 | | | | |
| 85.90 | 73 | | | | |
| 87.00 | 6138 | | | | |
| 88.00 | 6183 | | | | |
| 90.90 | 557 | | | | |
| 92.00 | 3897 | | | | |
| 93.00 | 5449 | | | | |
| 94.00 | 15249 | | | | |
| 95.00 | 139197 | | | | |
| 96.00 | 9622 | | | | |
| 97.10 | 320 | | | | |
| 103.90 | 653 | | | | |
| 104.90 | 196 | | | | |

| m/z | abund. | | | | |
|--------|--------|--|--|--|--|
| 118.90 | 674 | | | | |
| 127.95 | 479 | | | | |
| 128.90 | 136 | | | | |
| 129.85 | 426 | | | | |
| 130.90 | 130 | | | | |
| 134.90 | 300 | | | | |
| 136.90 | 242 | | | | |
| 140.95 | 1253 | | | | |
| 141.80 | 59 | | | | |
| 142.95 | 1244 | | | | |
| 145.90 | 63 | | | | |
| 147.00 | 50 | | | | |
| 147.95 | 302 | | | | |
| 154.95 | 289 | | | | |
| 156.95 | 171 | | | | |
| 158.80 | 59 | | | | |
| 160.80 | 58 | | | | |
| 161.00 | 63 | | | | |
| 171.70 | 60 | | | | |
| 173.90 | 98701 | | | | |
| 174.95 | 7288 | | | | |
| 175.90 | 96045 | | | | |

| | | _ | | | - | | |
|-------|-------|---|--------|-----|---|--------|------|
| 64.05 | 470 | | 105.95 | 611 | | 176.90 | 6065 |
| 67.05 | 432 | | 110.80 | 64 | | 177.95 | 142 |
| 68.00 | 15313 | | 114.90 | 149 | | 207.95 | 129 |
| 69.00 | 14886 | | 115.95 | 474 | | 253.00 | 29 |
| 70.05 | 1169 | | 116.90 | 778 | | | |
| 72.05 | 650 | | 117.90 | 474 | | | |
| | • | • | | • | • | - | |

| Mass | Criteria | % Rel. Abundance |
|------|---------------------------------------|------------------|
| 50 | 15-40% of mass 95 | 22.20% |
| 75 | 30-80% of mass 95 | 51.48% |
| 95 | Base peak, 100% of relative abundance | 100.00% |
| 96 | 5-9% of mass 95 | 6.91% |
| 173 | <2% of mass 174 | 0.00% |
| 174 | 50-120% of mass 95 | 70.91% |
| 175 | 5-9% of mass 174 | 7.38% |
| 176 | 95-101% of mass 174 | 97.31% |
| 177 | 5-9% of mass 176 | 6.31% |

Initial Calibration

Initial calibration was performed for both soil and water analyses. Both calibration curves included at least five calibration points, were performed prior to ICVs, samples and blanks, and were within twelve hours of a BFB check standard. Relative Response Factors (RRF) and Percent Relative Standard Deviations (%RSD) were calculated for each analyte and surrogate. The mean RRF and %RSD for target analytes were within control limits as stated in Table 18 of the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017. Below, the RRF, mean RRF and %RSDs were recalculated from one target analyte and surrogate associated with each internal standard and compared to laboratory reported values for the water calibration.

Water Calibration

| Internal Standard | Target Analyte | Calibration Level | IS Response | Analyte Response | RRF | Mean RRF | %RSD |
|-------------------------|----------------------|----------------------|----------------|---------------------|-------|-------------|--------|
| | | 2 | 268150 | 3752 | 0.35 | 0.509 | 15.837 |
| | | 5 | 264492 | 12849 | 0.486 | | |
| | Dichloro- | 10 | 259432 | 25657 | 0.494 | | |
| Pentafluoro- benzene | difluoro- methane | 20 | 271837 | 57930 | 0.533 | | |
| Denzerie | | 50 | 270155 | 151172 | 0.56 | | |
| | | 100 | 268034 | 323388 | 0.603 | | |
| | | 200 | 289910 | 625892 | 0.54 | | |

| | | 0.2 | 286150 | 103161 | 0.361 | 0.365 | 0.766 |
|-----------------------|------------------------|-----|--------|--------|-------|-------|--------|
| | - | 0.5 | 276425 | 100477 | 0.363 | | |
| | | 1 | 273251 | 99741 | 0.365 | | |
| | | 2 | 268150 | 98350 | 0.367 | | |
| | Dibromo- | 5 | 264492 | 96517 | 0.365 | | |
| | fluoromethane (s) | 10 | 259432 | 94385 | 0.364 | | |
| | (5) | 20 | 271837 | 99381 | 0.366 | | |
| | | 50 | 270155 | 98231 | 0.364 | | |
| | | 100 | 268034 | 97945 | 0.365 | | |
| | | 200 | 289910 | 107672 | 0.371 | | |
| | | 0.5 | 292406 | 1167 | 0.399 | 0.524 | 14.623 |
| | | 1 | 289379 | 3401 | 0.588 | | |
| | | 2 | 284987 | 4391 | 0.385 | | |
| | | 5 | 280557 | 15399 | 0.549 | | |
| | Carbon | 10 | 274828 | 29030 | 0.528 | | |
| | Tetrachloride | 20 | 293198 | 66222 | 0.565 | | |
| | | 50 | 290253 | 163881 | 0.565 | | |
| | | 100 | 289787 | 335583 | 0.579 | | |
| | | 200 | 319832 | 711557 | 0.556 | | |
| 1,4-difluoro- | | 0.2 | 300060 | 335276 | 1.117 | 1.113 | 0.716 |
| benzene | - | 0.5 | 292406 | 325733 | 1.114 | | |
| | | 1 | 289379 | 320340 | 1.107 | | |
| | | 2 | 284987 | 315396 | 1.107 | | |
| | | 5 | 280557 | 312969 | 1.116 | | |
| | Toluene-d8 (s) | 10 | 274828 | 306258 | 1.114 | | |
| | | 20 | 293198 | 321156 | 1.095 | | |
| | | 50 | 290253 | 323162 | 1.113 | | |
| | | 100 | 289787 | 325490 | 1.123 | | |
| | | 200 | 319832 | 358301 | 1.12 | | |
| | | 0.5 | 231509 | 726 | 0.314 | 0.392 | 12.774 |
| | | 1 | 229196 | 2116 | 0.462 | | |
| | | 2 | 222484 | 2741 | 0.308 | | |
| | | 5 | 221240 | 9223 | 0.417 | | |
| Chloro- benzene-d5 | Tetrachloro- ethene | 10 | 216758 | 17051 | 0.393 | | |
| Delizerie-us | | 20 | 228386 | 38209 | 0.418 | | |
| | | 50 | 231775 | 94064 | 0.406 | | |
| | [| 100 | 232198 | 192216 | 0.414 | | |
| | [| 200 | 254687 | 408369 | 0.401 | | |
| | Isopropyl- | 0.2 | 127486 | 2127 | 4.171 | 3.478 | 12.094 |
| | benzene | 0.5 | 123536 | 3619 | 2.93 | | |

| 1 | | | | | | | |
|-------------------|------------------------|-----|--------|---------|-------|-------|------|
| | | 1 | 122860 | 9770 | 3.976 | | |
| | | 2 | 119482 | 13197 | 2.761 | | |
| | | 5 | 119248 | 42286 | 3.546 | | |
| | | 10 | 117909 | 77767 | 3.298 | | |
| | | 20 | 127001 | 177377 | 3.492 | | |
| | | 50 | 127415 | 446470 | 3.504 | | |
| | | 100 | 129323 | 915600 | 3.54 | | |
| | | 200 | 136940 | 1954165 | 3.568 | | |
| 1,4- dichloro- | | 0.2 | 127486 | 116647 | 0.915 | 0.921 | 1.08 |
| benzene-d4 | | 0.5 | 123536 | 115532 | 0.935 | | |
| | | 1 | 122860 | 113267 | 0.922 | | |
| | | 2 | 119482 | 111235 | 0.931 | | |
| | 4-bromo- | 5 | 119248 | 110383 | 0.926 | | |
| | fluoro- benzene (s) | 10 | 117909 | 108881 | 0.923 | | |
| | 001120110(0) | 20 | 127001 | 114769 | 0.904 | | |
| | | 50 | 127415 | 116941 | 0.918 | | |
| | | 100 | 129323 | 117667 | 0.91 | | |
| | | 200 | 136940 | 127444 | 0.931 | | |

Initial Calibration Verification

A second source calibration verification was analyzed after the soil and water calibrations. The RRF for target analytes and surrogate analytes were above minimums presented in the CLP Guidance. The percent difference between the mean RRF of the initial calibration and the initial calibration verification were within control limits stated by the CLP Guidance.

Continuing Calibration Verification

Continuing calibration verifications were evaluated to determine instrument performance and calibration validity during each analytical run. The RRF %D was outside of the control limit for carbon tetrachloride, trichlorofluoromethane and 1,1,1-trichloroethane, associated with water sample data. Data for the Trip and Field blanks are considered estimated for carbon tetrachloride. The CCV for soil data had 1,1-dichloroethene and trans-1,2-dichloroethene outside of control limits for %D indicated in Table 18. These analytes are UJ flagged in the data table. The RRF of analytes for CCVs was within control limits.

Blanks

Method blanks, field blanks and trip blanks were analyzed to assess possible contamination from the laboratory or the field. One method blank was analyzed per analytical batch. One trip blank and one field blank were provided per SDG. There were no detections in any blanks associated with this SDG.

Surrogates

The percent recovery of surrogate compounds were evaluated to determine method and instrument performance.

Surrogate recoveries for water samples, soil samples and quality control samples were within laboratory control limits. Raw data was checked for calculation and transcription errors and none were found.

Laboratory Control Sample

One laboratory control sample was analyzed per analytical batch included in the SDG. The percent recoveries for target analytes were within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

One matrix spike was analyzed for the analytical batch encompassing soil data within the report. Percent recoveries for this matrix spike were within laboratory and CLP control limits.

A matrix spike and matrix spike duplicate were analyzed for the analytical batch encompassing water data within the report. The percent recovery for benzene was below the lower control limit. Because the original sample used as the source for the MS and MSD was not included in the SDG and the LCS percent recovery for benzene was within control limits, no data was qualified in the data table.

Laboratory Duplicate

A laboratory duplicate was performed on sample JC12861-1 (MW-1I-20'). RPD values were outside of laboratory control limits for all detected analytes. Only analytes with concentrations greater than 5 times the reporting limit are controlled, which was none of the analytes detected in the original sample and the duplicate.

Internal Standard

All required internal standards were added to samples, blank samples and quality control samples at the same specified concentration to monitor system performance and quantify target analytes. Raw data was checked against acceptance criteria for retention times and area response. Retention times did not differ more than 10 seconds from the check standard (beginning calibration check verification). Area responses were within 50-200% of the check standard. Below are the area responses of internal standards for specific samples and check standards. Retention time is expressed in minutes.

| | Tert Butyl Alcohol-D9 | | | Pentafluorobenzene | | | | 1,4- | Difluorok | enzene | ; | |
|---|-----------------------|------------------------|-----|--------------------|--------|------------------------|-----------|---------|-----------|------------------------|-----------|---------|
| | Area | % of Check Stnd. | RT | ∆ RT | Area | % of Check Stnd. | RT | ∆ RT | Area | % of Check Stnd. | RT | ∆ RT |
| Check Standard (1/21/16 21:51) | 72158 | 100% | 8.3 | 0 | 239231 | 100% | 10.7 5 | 0 | 250589 | 100% | 11.6 8 | 0 |
| V2A7019- MB1 | 80223 | 111% | 8.3 | 0.0 1 | 251717 | 105% | 10.7 5 | 0 | 260491 | 104% | 11.6 8 | 0 |
| V2A7019- BS | 71606 | 99% | 8.3 | 0 | 235863 | 99% | 10.7 5 | 0 | 246965 | 99% | 11.6 8 | 0 |
| JC12916- 14MS | 71288 | 99% | 8.3 | 0 | 235177 | 98% | 10.7 5 | 0 | 250884 | 100% | 11.6 8 | 0 |
| JC12916- 14MSD | 78946 | 109% | 8.3 | 0.0 1 | 256583 | 107% | 10.7 5 | 0 | 273611 | 109% | 11.6 8 | 0 |
| JC12861- 5 | 84942 | 118% | 8.3 | 0.0 1 | 252726 | 106% | 10.7 5 | 0 | 261828 | 104% | 11.6 8 | 0 |
| JC12861- 6 | 83683 | 116% | 8.3 | 0.0 1 | 246699 | 103% | 10.7 5 | 0 | 251323 | 100% | 11.6 8 | 0 |

| | | Chlorobe | nzene-D5 | | 1,4-Dichlorobenzene-d4 | | | |
|--------------------------------------|--------|------------------------|----------|-----|------------------------|------------------------|-------|-----|
| | Area | % of Check Stnd. | RT | ΔRT | Area | % of Check Stnd. | RT | ΔRT |
| Check Standard (1/21/16 21:51) | 202912 | 100% | 14.67 | 0 | 113053 | 100% | 16.83 | 0 |
| V2A7019-MB1 | 206171 | 102% | 14.67 | 0 | 110690 | 98% | 16.83 | 0 |
| V2A7019-BS | 199816 | 98% | 14.67 | 0 | 112664 | 100% | 16.83 | 0 |
| JC12916-14MS | 202176 | 100% | 14.67 | 0 | 115159 | 102% | 16.83 | 0 |
| JC12916- 14MSD | 217501 | 107% | 14.67 | 0 | 118415 | 105% | 16.83 | 0 |
| JC12861-5 | 207241 | 102% | 14.67 | 0 | 112126 | 99% | 16.83 | 0 |
| JC12861-6 | 199811 | 98% | 14.67 | 0 | 109434 | 97% | 16.83 | 0 |

| | Tert | Butyl Alc | ohol-D | 9 | Pen | tafluorob | enzene | ż | 1,4- | Difluoro | penzene | 5 |
|---|--------|------------------------|--------|----------|--------|------------------------|--------|----------|--------|-----------------------|-----------|---------|
| | Area | % of Check Stnd. | RT | ∆ RT | Area | % of Check Stnd. | RT | ∆ RT | Area | % of Check Stnd | RT | Δ RT |
| Check Standard (1/24/16 11:12) | 189750 | 100% | 7.25 | 0 | 420164 | 100% | 9.46 | 0 | 571409 | 100% | 10.3 7 | 0 |
| VY7020- MB | 165361 | 87% | 7.25 | 0 | 401290 | 96% | 9.45 | 0.0 1 | 543143 | 95% | 10.3 7 | 0 |
| VY7020- BS | 193132 | 102% | 7.26 | 0.0 1 | 413719 | 98% | 9.45 | 0.0 1 | 568506 | 99% | 10.3 7 | 0 |
| JC12861- 1 | 214359 | 113% | 7.25 | 0 | 427363 | 102% | 9.45 | 0.0 1 | 588133 | 103% | 10.3 7 | 0 |
| JC12861- 2 | 189375 | 100% | 7.25 | 0 | 439819 | 105% | 9.46 | 0 | 605136 | 106% | 10.3 7 | 0 |
| JC12861- 3 | 201423 | 106% | 7.25 | 0 | 445684 | 106% | 9.46 | 0 | 606690 | 106% | 10.3 7 | 0 |
| JC12861- 8 | 180821 | 95% | 7.24 | 0.0 1 | 441306 | 105% | 9.45 | 0.0 1 | 599111 | 105% | 10.3 7 | 0 |
| JC12861- 1DUP | 213763 | 113% | 7.24 | 0.0 1 | 435556 | 104% | 9.45 | 0.0 1 | 602841 | 106% | 10.3 7 | 0 |
| JC12861- 2MS | 151785 | 80% | 7.25 | 0 | 445731 | 106% | 9.45 | 0.0 1 | 605113 | 106% | 10.3 7 | 0 |
| JC12861- 7 | 184415 | 97% | 7.25 | 0 | 432639 | 103% | 9.45 | 0.0 1 | 592423 | 104% | 10.3 7 | 0 |
| JC12861- 9 | 190434 | 100% | 7.25 | 0 | 425229 | 101% | 9.45 | 0.0 1 | 587888 | 103% | 10.3 7 | 0 |

| | Chlorobenzene-D5 | | | | | 1,4-Dichloro | benzene-d4 | ŀ |
|---|------------------|------------------------|-------|-----|--------|------------------------|------------|-----|
| | Area | % of Check Stnd. | RT | ΔRT | Area | % of Check Stnd. | RT | ΔRT |
| Check Standard (1/24/16 11:12) | 446881 | 100% | 13.52 | 0 | 214290 | 100% | 15.82 | 0 |
| VY7020-MB | 432731 | 97% | 13.52 | 0 | 198556 | 93% | 15.82 | 0 |
| VY7020-BS | 440060 | 98% | 13.52 | 0 | 214777 | 100% | 15.82 | 0 |
| JC12861-1 | 470646 | 105% | 13.52 | 0 | 210797 | 98% | 15.82 | 0 |
| JC12861-2 | 475030 | 106% | 13.52 | 0 | 212223 | 99% | 15.82 | 0 |
| JC12861-3 | 477033 | 107% | 13.52 | 0 | 210846 | 98% | 15.82 | 0 |
| JC12861-8 | 466068 | 104% | 13.52 | 0 | 208497 | 97% | 15.82 | 0 |
| JC12861- 1DUP | 474838 | 106% | 13.52 | 0 | 213056 | 99% | 15.82 | 0 |
| JC12861-2MS | 454710 | 102% | 13.52 | 0 | 209961 | 98% | 15.82 | 0 |
| JC12861-7 | 467252 | 105% | 13.52 | 0 | 207020 | 97% | 15.82 | 0 |
| JC12861-9 | 469844 | 105% | 13.52 | 0 | 208812 | 97% | 15.82 | 0 |

2.4.2 Percent Moisture

Percent moisture was performed and calculated on two soil samples for SDG JC12861. Samples results were verified by recalculation of raw data, where; *Percent Moisture* = $\frac{Wet Weight-Dry Weight}{Wet Weight-Tare Weight} x100\%$.

2.5 SDG JC14061

Data validation and review was conducted for soil samples within SDG JC14061 in accordance with the EPA National Functional Guidelines for Superfund Organic Methods Review, January 2017, EPA National Functional Guidelines for Superfund Inorganic Methods Review, January 2017 and NYDEC Analytical Services Protocol. The SDG included seven groundwater samples and one field blank to be analyzed for VOCs by EPA 624, SVOCs by EPA 625, pesticides and PCBs by EPA 608, herbicides by EPA 8151, metals by EPA 6010C and mercury by EPA 7470A. Samples specific to JC14061 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|-----------|-------------|-------------|-------------|
| JC14061-1 | MW-1I | Groundwater | 2/10/2016 | 13:13 |
| JC14061-2 | MW-1S | Groundwater | 2/10/2016 | 14:14 |
| JC14061-3 | MW-2 | Groundwater | 2/10/2016 | 11:08 |
| JC14061-4 | MW-3 | Groundwater | 2/10/2016 | 16:13 |
| JC14061-5 | MW-4S | Groundwater | 2/10/2016 | 13:35 |
| JC14061-6 | MW-4I | Groundwater | 2/10/2016 | 14:45 |
| JC14061-7 | Field/FB | Groundwater | 2/10/2016 | 11:16 |
| JC14061-8 | MW-A | Groundwater | 2/10/2016 | 12:00 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey the same day as the samples were collected. As indicated by the analytical laboratory sample receipt summary: sample coolers were received with temperatures below 6°C, VOAs to be analyzed for VOCs did not contain headspace, sufficient sample volume was received, custody seals were intact and containers were accurately labeled and reflected the chain of custody.

2.5.1 VOCs

Preservation and Holding Time

Samples were received within the required temperature range and were analyzed for VOCs between one and two days after collection. VOAs were not preserved.

GC/MS Instrument Performance Check

Ion abundances were averaged over three scans and background was subtracted from less than 20 scans away. BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence, per EPA method 624.

Initial Calibration

RRF and %RSD were within control limits for all calibrations associated with SDG JC14061. A minimum of 5 calibration points was used for the calibration curves.

Initial Calibration Verification

An ICV was analyzed after the last ICAL standard. The ICV RRF for target analytes was greater than the minimum RRF. The %D between the ICAL mean RRF and the ICV RRF was within limits of the CLP guidance for trace volatiles.

Continuing Calibration Verification

Target analytes were within %D and RT limits for opening CCVs. RRFs were greater than minimums presented in Table 4 of the CLP guidance.

Blanks

No target analytes were detected in the field blank or method blanks.

Surrogates

Surrogate recoveries in environmental samples and quality control samples were within control limits.

Laboratory Control Samples

A LCS was included in both analytical batches with samples from SDG JC14061. The recovery of target analytes was within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

Matrix spike and matrix spike duplicates were analyzed from sources outside of SDG JC14061. Analytes outside of control limits are not expected to affect data quality of JC14061 samples since the LCS and laboratory duplicate were within control limits.

Laboratory Duplicate

A laboratory duplicate was analyzed as part of the laboratory quality control samples. The source of the laboratory duplicate was not part of SDG JC14061; all RPD values were within control limits.

Internal Standards

The retention time of the internal standards was compared against daily beginning continuing calibration verifications. The retention times did not differ more than 10 seconds and area counts were within 50-200% of the check standard.

Target Analyte Identification

Reported target analytes are accurately identified based on ion abundances, analyte spectrums and retention times.

2.5.2 SVOCs

Preservation and Holding Times

Samples to be prepared for SVOC analysis were extracted February 11, 2016 and analyzed February 12, 2016. Extraction and analysis dates were confirmed with preparation logs and instrument run logs. Samples were extracted and analyzed within holding times.

GC/MS Instrument Performance Check

A DFTPP standard was analyzed at the beginning of every analytical sequence that included analysis of calibration standards, samples associated with JC14601 and quality control samples. The last sample in the analytical sequences were analyzed within 12 hours of the DFTPP standard. Ion abundances were within CLP Guidance limits.

Initial Calibration

Target analytes had mean RRFs above the minimum RRF values stated in CLP guidance. The %RSD between the RRFs were below maximum control limits, except for 2,4-dinitirophenol, which was calibrated with a quadratic regression with a coefficient of 0.999.

Initial Calibration Verification

RRF of target analytes in the ICV were above minimum RRF stated by CLP guidance. The %D between the ICAL average and the ICV RRF for 2-methylnaphthalene was above the upper control limit of +20%. Results for 2-methlynaphthalene were not reported by the laboratory. The RT and %D for the remaining target analytes were within control limits.

Continuing Calibration Verification

The % drift between the ICAL and the CCV standard for 2,4-dinitrophenol was outside of the ±20% limit. The affected sample, JC14061-3 (MW-2), has the result for 2,4-dinitrophenol UJ flagged in the table. The other target analytes have RRFs above CLP guidance and %D within control limits.

Blanks

One method blank was extracted with JC14061-3 (MW-2) and a field blank (JC14061-7) was analyzed for SVOCs. No analytes were detected in either sample. Sample chromatograms were examined for potential signs of contamination and none were found.

Surrogates

The recovery of surrogate compounds was within control limits. Raw data was compared to reported surrogate recoveries and no discrepancies were found.

Laboratory Control Samples

One laboratory control sample was included in the extraction batch with samples from SDG JC14061. Recovery of target analytes was within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

The source for the MS/MSD was JC14061-3 (MW-2). The recovery of benzidine was below the lower control limit for the MS resulting in a RPD violation. Since the recovery of benzidine was 2% for the MS and benzidine is ND in the samples, results in the data table are qualified as unusable (R).

Internal Standard

The internal standard area counts in samples and quality control samples were within 50-200% of the opening CCV area counts. The retention time of internal standards did not differ more than 30 seconds from the opening CCV target analyte retention times.

Target Analyte Identification

No target analytes were identified in MW-2 and the field blank sample. Chromatogram were reviewed to ensure all measurable analyte peaks were accounted for.

2.5.3 Pesticides

Preservation and Holding Time

Samples were extracted two days after collection and were analyzed the same day as extraction. Extraction and analysis were within holding times.

GC/ECD Instrument Performance Check

The breakdown of DDT and endrin was below 20%. The breakdown check was performed daily before ICAL and sample analysis.

Initial Calibration

More than five calibration levels were performed for target analyte. Toxaphene and chlordane were calibrated using a one-point calibration curve. Mean retention times were calculated using an initial calibration standard and two subsequent continuing calibration verification standards. Response factors were calculated based on the internal standard area counts. The %RSD between the RFs of the calibration levels was less than 25%.

Continuing Calibration Verification

Calibration was verified at the beginning and the end of every analytical sequence. Retention times of analytes were within retention time windows. The %D between the CCV and the ICAL was within a ±25% range except for analytes on the second column of the last CCV. Since there were no detection of pesticides within the sample and the first column was within control limits, no data was qualified.

Blanks

The method blank extracted with the samples of JC14061 and the field blank did not contain any target analyte pesticides. Raw data was verified that no contamination was present.

Surrogates

Recovery of surrogates were within 30-150%. Retention times of surrogates were verified against retention time windows.

Laboratory Control Sample

One LCS/LCSD pair was extracted with samples. The recovery of endrin aldehyde, gamma-BHC, endosulfan sulfate and endrin were above CLP control limits. Because the samples did not contain any pesticides, no data was qualified in the data table.

2.5.4 PCBS

Preservation and Holding Time

Samples were extracted two days after collection and were analyzed the same day as extraction. Extraction and analysis were within holding times.

Initial Calibration

The calibration factor for Aroclor 1016 and 1260 peaks were calculated from a six-level calibration. The %RSD between calibration factors was less than 20% for both columns. Retention time windows (±0.03) were calculated from the ICAL and two continuing calibration verifications.

Continuing Calibration Verification

ICVs and CCV were within the inclusive range of ±25% %D when compared to CF of target analytes from the ICAL. CCVs were analyzed every ten samples and at the beginning and end of each sequence.

Blanks

A method blank was extracted with MW-2 and the field blank. No PCBs were found in the method blank or field blank and verified against raw data chromatograms.

Surrogates

Surrogate recoveries were within 30-150% of true values. Surrogate retention times were verified and were within RT windows.

Matrix Spike/Matrix Spike Duplicate

MW-2 was used as the source for the MS/MSD. The recovery and RPD of Aroclors 1016 and 1260 were within CLP Guidance limits.

Laboratory Control Sample

The recovery of Aroclors 1016 and 1260 were with CLP Guidance limits.

2.5.5 Herbicides

Preservation and Holding time

Samples were extracted two days after collection and were analyzed 4 days after extraction. Extraction and analysis were within holding times.

Initial Calibration

The calibration factor for target analyte herbicides were calculated from a minimum of five calibration points. The %RSD between calibration factors was less than 20% for both columns. Retention time windows (±3*StdDev) were

calculated from the ICAL and two continuing calibration verifications.

Continuing Calibration Verification

ICVs and CCV were within the inclusive range of $\pm 25\%$ %D when compared to CF of target analytes from the ICAL. CCVs were analyzed every ten samples and at the beginning and end of each sequence.

Blanks

A method blank was extracted with MW-2 and the field blank. No target analyte herbicides were found in the method blank or field blank which was verified against raw data chromatograms.

Surrogates

Surrogate recoveries were within 39-159% of true values. Surrogate retention times were verified and were within RT windows.

Matrix Spike/Matrix Spike Duplicate

MW-2 was used as the source for the MS/MSD. The recovery and RPD of target analyte herbicides were within laboratory control limits.

Laboratory Control Sample

The recovery of target analyte herbicides were within laboratory control limits.

2.5.6 Metals

Preservation and Holding Time

The water samples were collected in bottles containing nitric acid (HNO₃) to a pH<2. This was verified by the lab upon digestion of the samples, which was completed two days after collection. Samples were analyzed the same day as digestion. This was confirmed with raw data.

Calibration

The instrument was calibrated with a 2-point calibration curve and internal standard evaluation. Linear regression was used with a 1/conc weighting. The linear calibration range and instrument detection limits were evaluated on 12/9/15.

ICV and CCV recoveries of target analytes were within control limits. CCVs were analyzed every ten samples and followed by a calibration blank.

Blanks

A CCB was analyzed every 10 samples following the CCV sample. No analytes were detected above the reporting limits in CCBs. A method blank was digested along with MW-2 and the field blank. No analytes above the reporting limit were detected. The field blank also did not have any analytes detected above the laboratory reporting limit.

Interference Check Standard

The ICSA and ICSAB analytes were within $\pm 20\%$ of true values. Analytes that were not spiked within the ICSA were within instrument control limits.

Laboratory Control Sample

A LCS for total and dissolved analytes was prepared with JC14061 samples. The recovery of target analytes were within the 70-130% control limits.

Matrix Spike/Matrix Spike Duplicate

A MS/MSD were prepared with samples from SDG JC14061. The matrix source for the MS/MSD was not a sample from JC14061. Percent recoveries and RPD values were within laboratory control limits.

Serial Dilution

One serial dilution was performed on a sample not included in SDG JC14061. The %D between the original sample and the serial dilution were outside of control limits due to low analyte concentration.

2.5.7 Mercury

Preservation and Holding Time

Samples were preserved with nitric acid and were pH checked by the laboratory before digestion to confirm a pH<2. Samples were digested two days after collection and analyzed the same day as digestion.

Calibration

The mercury instrumentation was calibrated with six standards that resulted in a correlation coefficient of 0.999. Initial and continuing calibration verification was with the 85-115% control limits. CCVs were analyzed every ten samples.

Blanks

CCBs were analyzed immediately after CCVs and were below the mercury reporting limit. One method blank was prepared with samples from JC14061 and mercury was not detected above the reporting limit.

Laboratory Control Sample

One LCS was digested with samples from SDG JC14061. The recovery of mercury was within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

A MS/MSD were digested with samples from SDG JC14061. The source of the MS/MSD was not part of SDG JC14061. The recovery and RPD of mercury in the MS/MSD was within laboratory control limits.

2.6 SDG JC15505

Data validation and review was conducted for air samples within SDG JC15505 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017, EPA Method TO-15 and NYDEC Analytical Services Protocol. The SDG included three air samples to be analyzed for VOCs by TO-15. Samples specific to JC15505 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|-----------|--------|-------------|-------------|
| JC15505-1 | SV-11 | Air | 3/4/2016 | 11:36 |
| JC15505-2 | IAQ-11 | Air | 3/4/2016 | 11:36 |
| JC15505-3 | IAQ-10 | Air | 3/4/2016 | 11:47 |
| JC15505-4 | SV-9 | Air | 3/4/2016 | 11:54 |
| JC15505-5 | IAQ-9 | Air | 3/4/2016 | 11:54 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey on March 5, 2016. As indicated by the analytical laboratory sample receipt summary: custody seals were intact and containers were accurately labeled and reflected the chain of custody. The air samples were collected in 6-liter canisters and no damage was present when received by the laboratory.

Preservation and Holding Time

The air samples were collected in cleaned metal canisters. They arrived at the analytical laboratory one day after sampling and were analyzed one March 12, 2016 and March 13, 2016. Raw data was verified to ensure analysis date reported is correct.

GC/MS Instrument Performance Check

BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence and 24 hours within the last samples injected, per method TO-15.

Initial Calibration

RRF and %RSD were within control limits for all calibrations associated with SDG JC15505. A minimum of 5 calibration points were used for the calibration curves.

Initial Calibration Verification

An ICV standard was analyzed after each ICAL. Minimum RRF was met for target analytes and %D between the mean ICAL RRF and the ICV RRF was within control limits.

Continuing Calibration Verification

Target analytes were within %D limits for opening CCVs.

Blanks

Method blanks were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. No analytes other than surrogates and internal standards were detected in method blanks. Raw data was verified that no other peaks were present in the chromatograms.

Surrogates

The surrogate was within laboratory control limits. Raw data for samples was checked to ensure that when dilutions were run, both recoveries were within control limits.

Laboratory Control Samples

LCS/LCSDs were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. Recoveries and RPD values of target analytes were within control limits.

Laboratory Duplicate

A laboratory duplicate was performed on samples not included in SDG JC15505. All RPD values were within control limits.

Summa Cleaning Certification

Summa's were cleaned and no residual VOCs were found.

Internal Standards

The retention time of the internal standards was compared against daily opening continuing calibration verifications. The retention times did not differ more than 0.33 minutes and area counts were within 40% of the check standard for samples associated with JC15505.

Target Analyte Identification

Raw data was verified to ensure that analyte identification was accurate by comparing ion abundances, analyte mass spectra and retention times to check standards. Some target analytes did not meet the abundance criteria of $\pm 20\%$ and were accepted based on analyte spectra and retention times.

2.7 SDG JC15742

Data validation and review was conducted for soil samples within SDG JC15742 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 and NYDEC Analytical Services Protocol. The SDG included eight soil samples to be analyzed for VOCs by EPA 8260C and percent moisture by SM2540. Samples specific to JC15742 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|-----------|--------|-------------|-------------|
| JC15742-1 | MW-7-23 | Soil | 3/8/2016 | 9:10 |
| JC15742-2 | MW-7-34 | Soil | 3/8/2016 | 9:35 |
| JC15742-3 | MW-6-23 | Soil | 3/8/2016 | 13:23 |
| JC15742-4 | MW-6-30 | Soil | 3/8/2016 | 13:45 |
| JC15742-5 | MW-5-22.5 | Soil | 3/8/2016 | 14:30 |
| JC15742-6 | MW-5-30 | Soil | 3/8/2016 | 14:45 |
| JC15742-7 | MW-5-35 | Soil | 3/8/2016 | 14:55 |
| JC15742-8 | MW-5 | Soil | 3/9/2016 | 13:40 |

Samples were received within twenty-four hours of sampling to the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: the cooler temperature upon receipt was <6°C, custody seals were intact, containers were accurately labeled and reflected the chain of custody. Soil samples were collected using field core sampling with enclosed storage containers (EnCore[®] sampler) in compliance with preparation method EPA 5035 and were prepared by the analytical laboratory within twenty-four hours of sample receipt.

2.7.1 VOCs

Holding Times

Samples were received below 6^oC and within method holding limits by the analytical laboratory. Soil samples were prepared by the laboratory within 24 hours of receipt. Holding times for soil samples analyzed for VOCs were met.

Instrument Performance Check

Bromofluorobenzene (BFB) was evaluated for mass spectra requirements and frequency to ensure adequate instrument performance. BFB was analyzed prior to any samples, calibration standard, calibration verification and quality control samples. All samples were run within 12 hours of initial BFB injection. Ion abundances within the BFB peaks were within criteria as established in CLP guidance. Ion abundances were averaged over three scans and were background from less than 20 scans away.

Initial Calibration

Calibration curves included at least five calibration points, were performed prior to ICVs, samples and blanks, and were within twelve hours of a BFB check standard. Relative Response Factors (RRF) and Percent Relative Standard Deviations (%RSD) were calculated for each analyte and surrogate. The mean RRF and %RSD for target analytes were within control limits as stated in Table 18 of the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017.

Initial Calibration Verification

The RRF for target analytes and surrogate analytes were above minimums presented in the CLP Guidance. The

percent difference between the mean RRF of the initial calibration and the initial calibration verification were within control limits stated by the CLP Guidance.

Continuing Calibration Verification

Continuing calibration verifications were evaluated to determine instrument performance and calibration validity during each analytical run. The CCV analyzed on 3/14/2016 had a %D value outside of control limits for bromomethane. Bromomethane is UJ qualified in the data table for MW-5-30 and MW-5-35.

Blanks

Method blanks were analyzed to assess possible contamination from the laboratory. One method blank was analyzed per analytical batch. There were no detections in any blanks associated with this SDG.

Surrogates

The percent recovery of surrogate compounds was evaluated to determine method and instrument performance. Surrogate recoveries for soil and quality control samples were within laboratory control limits.

Laboratory Control Sample

One laboratory control sample was analyzed per analytical batch included in the SDG. The percent recoveries for target analytes were within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

One matrix spike was analyzed per analytical batch included in the SDG. The sources of the MS/MSDS were from SDG 15742. The recovery of methyl acetate was outside of control limits for batch VI8494. It is not expected that results for associated samples in this batch are affected since the source was not part of SDG 15742 and the LCS had an acceptable recovery.

Laboratory Duplicate

A laboratory duplicate was performed on a sample not included in SDG JC15742. No analytes were detected in the source sample or the laboratory duplicate.

Internal Standard

All required internal standards were added to samples, blank samples and quality control samples at the same specified concentration to monitor system performance and quantify target analytes. Raw data was checked against acceptance criteria for retention times and area response. Retention times did not differ more than 10 seconds from the check standard (beginning calibration check verification). Area responses were within 50-200% of the check standard.

Target Analyte Identification

Target analyte detections are confirmed by ion fragments and retention times. Acetone had ion fragments outside of the acceptable abundance range as defined by the initial calibration. These results were accepted as positively

identified compounds based on RT and analyte mass spectra compared to the reference spectrum.

Target Analyte Quantitation and Reported Contract Required Quantitation Limits

Results for positively identified analytes and method reporting limits were calculated correctly by the laboratory and adjusted based on initial sample volumes and percent solids for soil samples. The percent solids in the soil samples was greater than 30%. Internal standards based on the ICAL were used for quantitation of analytes in samples and quality control samples. The same primary ion fragments were used as the ICAL and mean RRFs were the same.

2.7.2 Percent Moisture

Percent moisture was performed and calculated for soil samples in SDG JC15742. In addition, two soil samples from JC16088 were also used to calculate dry weight totals in JC15742. Samples results were verified by recalculation of raw data, where; $Percent \ Moisture = \frac{Wet \ Weight - Dry \ Weight}{Wet \ Weight - Tare \ Weight} x100\%$.

2.8 SDG JC15931

Data validation and review was conducted for air samples within SDG JC15931 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017, EPA method TO-15 and NYDEC Analytical Services Protocol. The SDG included three air samples to be analyzed for VOCs by TO-15. Samples specific to JC15931 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|-----------|--------|-------------|-------------|
| JC15931-1 | Ambient | Air | 3/9/2016 | 8:49 |
| JC15931-2 | SV-10 | Air | 3/9/2016 | 9:27 |
| JC15931-3 | IAQ-10 | Air | 3/9/2016 | 9:27 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: custody seals were intact and containers were accurately labeled and reflected the chain of custody. The air samples were collected in 6-liter canisters and no damage was present when received by the laboratory.

Preservation and Holding Time

The air samples were collected in cleaned metal canisters. They arrived at the analytical laboratory one day after sampling and were analyzed on 3/18/2016 and 3/19/2016. Samples were analyzed within method holding limits.

GC/MS Instrument Performance Check

BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence and within 24-hours of the final sample, per method TO-15. Ion abundances were averaged over three consecutive scans and no more than 20 scans away from the BFB peak.

Initial Calibration

RRF and %RSD were within control limits for all calibrations associated with SDG JC15931. A minimum of 5 calibration points were used for the calibration curves.

Initial Calibration Verification

An ICV standard was analyzed after each ICAL and before any samples or quality control samples. RRF values and %D were within control limits.

Continuing Calibration Verification

Target analytes were within %D limits for opening CCVs.

Blanks

Method blanks were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. No analytes other than surrogates and internal standards were detected in method blanks. Raw data was verified that no other peaks were present in the chromatograms.

Surrogates

The surrogate recoveries were within laboratory control limits.

Laboratory Control Samples

LCS/LCSDs were performed with each sequence where analysis of environmental samples or SCC samples were performed. Recoveries and RPD values of target analytes were within control limits except for 1,2,4-trichlorobenzene analyzed on 3/5/16. Because this LCS/LCSD was part of the QC for SCC samples, no data was qualified for sampled in this SDG.

Laboratory Duplicate

Laboratory duplicates were performed on samples not included in SDG JC15931. All RPD values were within control limits.

Summa Cleaning Certification

Summa's were cleaned and no residual VOCs were found.

Internal Standards

The retention time of the internal standards was compared against dialing continuing calibration verifications. The retention times did not differ more than 0.33 minutes and area counts were within 40% of the check standards.

Target Analyte Identification

Some target analytes did not meet the abundance criteria of $\pm 20\%$ and were accepted based on analyte spectra and retention times.

2.9 SDG JC16088

Data validation and review was conducted for soil and blank water samples within SDG JC16088 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 and NYDEC Analytical Services Protocol. The SDG included two soil samples, one trip blank and one field blank to be analyzed for VOCs by EPA 8260C and percent moisture by SM2540. The percent moisture was used to calculate dry-weight results for SDG JC15742. Trip and field blanks were sampled to show there was no contamination from transportation or field activities which includes samples from SDG JC15742. Samples specific to JC16088 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|-------------|--------|-------------|-------------|
| JC16088-1 | Trip Blank | Water | 3/10/2016 | 9:50 |
| JC16088-2 | Field Blank | Water | 3/10/2016 | 9:50 |
| JC16088-3 | MW-6 | Soil | 3/10/2016 | 10:50 |
| JC16088-4 | MW-7 | Soil | 3/10/2016 | 10:50 |

Samples were received within twenty-four hours of sampling to the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: the cooler temperatures upon receipt were $<6^{\circ}$ C, custody seals were intact, containers were accurately labeled and reflected the chain of custody and there was no headspace in VOA (volatile organic analysis) containers. The QC blank samples were preserved with hydrochloric acid (HCl) and pH checked by the laboratory.

2.9.1 VOCs

Holding Times

QC blanks samples were analyzed 5 days after collection and within hold time.

Instrument Performance Check

Bromofluorobenzene (BFB) was evaluated for mass spectra requirements and frequency to ensure adequate instrument performance. BFB standards were analyzed at the beginning of every sequence and within 12 hours of the final sample injected. Ion abundances were within control limits and were averaged over three scans within the BFB peak and background subtracted from less than 20 scans away.

Initial Calibration

The calibration included at least five calibration points, was performed prior to ICVs, samples and blanks, and was within twelve hours of a BFB check standard. Relative Response Factors (RRF) and Percent Relative Standard Deviations (%RSD) were calculated for each analyte and surrogate. The mean RRF and %RSD for target analytes were within control limits.

Initial Calibration Verification

A second source calibration verification was analyzed after the soil and water calibrations. The percent difference

between the mean RRF of the initial calibration and the initial calibration verification were within control limits stated by the CLP Guidance.

Continuing Calibration Verification

Continuing calibration verifications were evaluated to determine instrument performance and calibration validity during each analytical run. The percent difference between the mean RRF of the initial calibration and the CCV were within control limits stated by the CLP Guidance.

Blanks

Method blanks, field blanks and trip blanks were analyzed to assess possible contamination from the laboratory or the field. One method blank was analyzed per analytical batch. There were no detections in any blanks associated with this SDG.

Surrogates

Surrogate recoveries were within control limits.

Laboratory Control Sample

One laboratory control sample was analyzed per analytical batch included in the SDG. The percent recoveries for target analytes were within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

A matrix spike was analyzed with the blank QC samples. Target analytes were within control limits.

Laboratory Duplicate

A laboratory duplicate was analyzed with the blank QC samples. RPD values were within laboratory control limits.

Internal Standard

All required internal standards were added to samples, blank samples and quality control samples at the same specified concentration to monitor system performance and quantify target analytes. Raw data was checked against acceptance criteria for retention times and area response. Retention times did not differ more than 10 seconds from the check standard (beginning calibration check verification). Area responses were within 50-200% of the check standard

2.9.2 Percent Moisture

Percent moisture was performed and calculated on two soil samples for SDG JC16088. Samples results were verified by recalculation of raw data, where; *Percent Moisture* = $\frac{Wet Weight-Dry weight}{Wet Weight-Tare Weight} x100\%$.

2.10 SDG JC16571

Data validation and review was conducted for air samples within SDG JC16571 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017, EPA Method TO-15 and NYDEC Analytical Services Protocol. The SDG included four air samples to be analyzed for VOCs by TO-15. Samples specific to JC16571 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|----------------|--------|-------------|-------------|
| JC16571-1 | SV-14 | Air | 3/17/2016 | 7:48 |
| JC16571-2 | SV-13 | Air | 3/17/2016 | 7:52 |
| JC16571-3 | SV-12 | Air | 3/17/2016 | 7:58 |
| JC16571-4 | Ambient Upwind | Air | 3/17/2016 | 8:02 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey on March 18, 2016. As indicated by the analytical laboratory sample receipt summary: custody seals were intact and containers were accurately labeled and reflected the chain of custody. The air samples were collected in 6-liter canisters and no damage was present when received by the laboratory.

Preservation and Holding Time

The air samples were collected in cleaned metal canisters and were analyzed within method holding limits. They arrived at the analytical laboratory one day after sampling and were analyzed between March 24, 2016 and March 26, 2016. Raw data was verified to ensure analysis dates reported are correct.

GC/MS Instrument Performance Check

BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence and 24 hours within the last samples injected, per method TO-15. Ion abundances were averaged over three scans and were background subtracted from less than 20 scans away.

Initial Calibration

RRF and %RSD were within control limits for all calibrations associated with SDG JC16571. A minimum of 5 calibration points were used for the calibration curves. The %RSD for tetrachloroethene on 3/1/2016 was outside of control limits. No data was qualified since the affected samples were SCCs.

Initial Calibration Verification

An ICV standard was analyzed after each ICAL. Minimum RRF was met for target analytes and %D between the mean ICAL RRF and the ICV RRF was within control limits.

Continuing Calibration Verification

Target analytes were within %D limits for opening CCVs.

Blanks

Method blanks were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. No analytes other than surrogates and internal standards were detected in method blanks. Raw data was verified that no other peaks were present in the chromatograms.

Surrogates

The recovery of the surrogate compound was within control limits for samples and associated quality control samples within the SDG.

Laboratory Control Samples

LCS/LCSDs were analyzed with each sequence where analysis of environmental samples or SCC samples were performed. Recoveries and RPD values of target analytes were within control limits.

Laboratory Duplicate

The source for the laboratory duplicate analyzed was not part of SDG JC16571. All RPD values were <20%.

Summa Cleaning Certification

Summa's were cleaned and no residual VOCs were found.

Internal Standards

The retention time of the internal standards was compared against daily opening continuing calibration verifications. The retention times did not differ more than 0.33 minutes and area counts were within 40% of the check standard.

Target Analyte Identification

Raw data was verified to ensure that analyte identification was accurate by comparing ion abundances, analyte mass spectra and retention times to check standards. Some target analytes did not meet the abundance criteria of $\pm 20\%$ due to matrix interferences but RT and mass spectra suggest correct analyte identification.

2.11 SDG JC16575

Data validation and review was conducted for soil samples within SDG JC16575 in accordance with the EPA National Functional Guidelines for Superfund Organic Methods Review, January 2017, EPA National Functional Guidelines for Superfund Inorganic Methods Review, January 2017, and NYDEC Analytical Services Protocol. The SDG included three groundwater samples and two quality control blank samples to be analyzed for VOCs by EPA 624, SVOCs by EPA 625, chlorinated herbicides by EPA 8151, chlorinated pesticides and PCBs by EPA 608, metals by EPA 6010C and mercury by EPA 7470A. Samples specific to JC16575 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|-----------|-------------|----------------|-------------|
| JC16575-1 | MW-5 | Groundwater | 3/18/2016 | 11:20 |
| JC16575-2 | MW-6 | Groundwater | 3/18/2016 | 11:42 |
| JC16575-3 | MW-7 | Groundwater | 3/18/2016 | 10:20 |
| JC16575-4 | FB | Water | 3/18/2016 | 9:10 |
| JC16575-5 | ТВ | Water | 3/15/2016 | 9:00 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey the same day as the samples were collected. As indicated by the analytical laboratory sample receipt summary: sample coolers were received with temperatures below 6°C, VOAs to be analyzed for VOCs did not contain headspace, sufficient sample volume was received, custody seals were intact and containers were accurately labeled and reflected the chain of custody.

2.11.1 VOCs

Preservation and Holding Time

Samples were received within the required temperature range and were analyzed for VOCs day after collection. VOAs were not preserved.

GC/MS Instrument Performance Check

Ion abundances were averaged over three scans and background was subtracted from less than 20 scans away. BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence and within 12-hours of the final sample injection time.

Initial Calibration

RRF and %RSD were within control limits. A minimum of 5 calibration points was used for the calibration curves.

Initial Calibration Verification

An ICV was analyzed after the last ICAL standard. The ICV RRF for target analytes was greater than the minimum RRF. The %D between the ICAL mean RRF and the ICV RRF was within limits of the CLP guidance for trace volatiles.

Continuing Calibration Verification

Target analytes were within %D and RT limits for opening CCVs. RRFs were greater than minimums presented in Table 4 of the CLP guidance.

Blanks

No target analytes were detected in the field blank, trip blank or method blanks. Contamination was not present in

these blanks.

Surrogates

Surrogate recoveries in environmental samples and quality control samples were within control limits.

Laboratory Control Samples

A LCS was included in both analytical batches with samples from SDG JC16575. The recovery of target analytes was within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

One matrix spike was analyzed from JC16575-2 (MW-6). The recovery of target analytes was within control limits.

Laboratory Duplicate

A laboratory duplicate was analyzed as part of the laboratory quality control samples. The source of the laboratory duplicate was JC16575-3 (MW-7). The target analyte cis-1,2-dichloroethene was outside of the RPD control limit. Sample data was accepted since the concentration of cis-1,2-dichloroethene was less than the reporting limit and cannot be controlled.

Internal Standards

The retention time of the internal standards was compared against daily beginning continuing calibration verifications. The retention times did not differ more than 10 seconds and area counts were within 50-200% of the check standard.

Target Analyte Identification

Reported target analytes are accurately identified based on ion abundances, analyte spectrums and retention times.

2.11.2 SVOCs

Preservation and Holding Times

Samples to be prepared for SVOC analysis were extracted three days after collected and analyzed one day after extraction. Samples were prepared and analyzed within holding time.

GC/MS Instrument Performance Check

A DFTPP standard was analyzed at the beginning of every analytical sequence that included analysis of calibration standards, samples associated with JC16575 and quality control samples. Per EPA 625, DFTPP was analyzed at the beginning of the day.

Initial Calibration

Target analytes had mean RRFs above the minimum RRF values stated in CLP guidance. The %RSD between the RRFs

were above maximum control limits, except for 2,4-dinitirophenol and 4,6-dinitro-2-methylphenol, which were calibrated with quadratic regressions that had coefficients of 0.999.

Initial Calibration Verification

RRF of target analytes in the ICV were above minimum RRF stated by CLP guidance. The %D between the ICAL average and the ICV RRF for 4-chloro-3-methylphenol was above the upper control limit of +20%. Results for 4-chloro-3-methylphenol were qualified with a UJ in the data table as estimated values. The RT and %D for the remaining target analytes were within control limits.

Continuing Calibration Verification

Analytes reported by the laboratory had RRF values above minimums and %RSD values between the ICAL mean and the CCV within recommended by CLP guidance and %RSD values

Blanks

One method blank was extracted with JC16575-1 (MW-5) and a field blank (JC16575-4) was analyzed for SVOCs. No analytes were detected in either sample. Sample chromatograms were examined for potential signs of contamination and none were found.

Surrogates

The recovery of surrogate compounds was within control limits. Raw data was compared to reported surrogate recoveries and no discrepancies were found.

Laboratory Control Samples

One laboratory control sample was included in the extraction batch with samples from SDG JC16575. Recovery of target analytes was within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

The source for the MS/MSD was JC16575-1 (MW-5). The recovery of benzidine was below the lower control limit for the MS. Since the recovery of benzidine was 9% for the MS and benzidine is ND in the samples, results in the data table are qualified as unusable (R).

Internal Standard

The internal standard area counts in samples and quality control samples were within 50-200% of the opening CCV area counts. The retention time of internal standards did not differ more than 30 seconds from the opening CCV target analyte retention times.

Target Analyte Identification

No target analytes were identified in MW-5 and the field blank sample. Chromatograms were reviewed to ensure all measurable analyte peaks were accounted for.

2.11.3 Pesticides

Preservation and Holding Time

Samples were extracted four days after collection and were analyzed two days after extraction. Extraction and analysis were within holding times.

GC/ECD Instrument Performance Check

The breakdown of DDT and endrin was below 20%. The breakdown check was performed before ICAL and prior to sample analysis.

Initial Calibration

More than five calibration levels were performed for target analyte. Toxaphene and chlordane were calibrated using a one-point calibration curve. Mean retention times were calculated using an initial calibration standard and two subsequent continuing calibration verification standards. Response factors were calculated based on the internal standard area counts. The %RSD between the RFs of the calibration levels were less than 25%.

Continuing Calibration Verification

Calibration was verified at the beginning and the end of every analytical sequence. Retention times of analytes were within retention time windows. The %D between the CCV and the ICAL was within a ±25%.

Blanks

The method blank extracted with the samples of JC16575 and the field blank did not contain any target analyte pesticides. Raw data was verified that no contamination was present.

Surrogates

Recovery of surrogates were within 30-150%. Retention times of surrogates were verified against retention time windows.

Matrix Spike/Matrix Spike Duplicate

The RPD value for endrin aldehyde, heptachlor and Aldrin were outside of CLP guidance or laboratory control limits. No data was qualified since the source sample was not part of SDG JC16575.

Laboratory Control Sample

One LCS was extracted with samples. The recovery of target analytes were within control limits.

2.11.4 PCBS

Preservation and Holding Time

Samples were extracted four days after collection and were analyzed two days after extraction. Extraction and analysis were within holding times.

Initial Calibration

The calibration factor for Aroclor 1016 and 1260 peaks were calculated from a six-level calibration. The %RSD between calibration factors was less than 20% for both columns. Retention time windows (±0.03) were calculated

from the ICAL and two continuing calibration verifications.

Continuing Calibration Verification

ICVs and CCV were within the inclusive range of ±25% %D when compared to CF of target analytes from the ICAL. CCVs were analyzed every ten samples and at the beginning and end of each sequence.

Blanks

A method blank was extracted with MW-5 and the field blank. No PCBs were found in the method blank or field blank and verified against raw data chromatograms.

Surrogates

Surrogate recoveries were within 30-150% of true values. Surrogate retention times were verified and were within RT windows.

Matrix Spike/Matrix Spike Duplicate

MW-5 was used as the source for the MS/MSD. The recovery and RPD of Aroclors 1016 and 1260 were within CLP guidance limits.

Laboratory Control Sample

The recovery of Aroclors 1016 and 1260 were with CLP guidance limits.

2.11.5 Herbicides

Preservation and Holding time

Samples were extracted four days after collection and were analyzed six days after extraction. Extraction and analysis were within holding times.

Initial Calibration

The calibration factor for target analyte herbicides were calculated from a minimum of five calibration points. The %RSD between calibration factors was less than 20% for both columns. Retention time windows (±3*StdDev) were calculated from the ICAL and two continuing calibration verifications.

Continuing Calibration Verification

ICVs and CCV were within the inclusive range of ±25% %D when compared to CF of target analytes from the ICAL. CCVs were analyzed every ten samples and at the beginning and end of each sequence.

Blanks

A method blank was extracted with MW-5 and the field blank. No target analyte herbicides were found in the method blank or field blank which was verified against raw data chromatograms.

Surrogates

Surrogate recoveries were within 39-159% of true values. Surrogate retention times were verified and were within RT windows.

Matrix Spike/Matrix Spike Duplicate

MW-5 was used as the source for the MS/MSD. The recovery and RPD of target analyte herbicides were within laboratory control limits.

Laboratory Control Sample

The recovery of target analyte herbicides were within laboratory control limits.

2.11.6 Metals

Preservation and Holding Time

The water samples were collected in bottles containing nitric acid (HNO_3). This was verified by the lab upon digestion of the samples. Samples were digested and analyzed within holding time.

Calibration

The instrument was calibrated with a 2-point calibration curve and internal standard evaluation. Linear regression was used with a 1/conc weighting. The linear calibration range and instrument detection limits were evaluated on 12/9/2015 and 1/26/2016, respectively.

ICV and CCV recoveries of target analytes were within control limits. CCVs were analyzed every ten samples followed by a calibration blank.

Blanks

A CCB was analyzed every 10 samples following the CCV sample. No analytes were detected above the reporting limits in CCBs. A method blank was digested along with MW-5 and the field blank. No analytes above the reporting limit were detected. The field blank also did not have any analytes detected above the laboratory reporting limit.

Interference Check Standard

The ICS analytes were within $\pm 20\%$ of true values. Analytes that were not spiked within the ICSA were within instrument control limits.

Laboratory Control Sample

A LCS for target analytes was prepared with JC16575 samples. The recovery of target analytes were within the 70-130% control limits.

Matrix Spike/Matrix Spike Duplicate

A MS/MSD were prepared with samples from SDG JC16575. The matrix source for the MS/MSD was not a sample from JC16575. Percent recoveries and RPD values were within laboratory control limits.

Serial Dilution

One serial dilution was performed on a sample not included in SDG JC16575. The %D between the original sample and the serial dilution were outside of control limits due to low analyte concentration.

2.11.7 Mercury

Preservation and Holding Time

Samples were preserved with nitric acid and were pH checked by the laboratory before digestion. Samples were digested four days after collection; within holding time limits.

Calibration

The mercury instrumentation was calibrated with six standards that resulted in a correlation coefficient of 0.999. Initial and continuing calibration verification was with the 85-115% control limits. CCVs were analyzed every ten samples.

Blanks

CCBs were analyzed immediately after CCVs and were below the mercury reporting limit. One method blank was prepared with samples from JC16575 and mercury was not detected above the reporting limit. Mercury was detected in the field blank. No data was qualified since MW-5 did not have a detectable concentration of mercury.

Laboratory Control Sample

One LCS was digested with samples from SDG JC16575. The recovery of mercury was within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

A MS/MSD were digested with samples from SDG JC16575. The source of the MS/MSD was not part of SDG JC16575. The recovery and RPD of mercury in the MS/MSD was within laboratory control limits.

2.12 SDG JC17397

Data validation and review was conducted for groundwater samples within SDG JC17397 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 and NYDEC Analytical Services Protocol. The SDG included one groundwater sample, one trip blank and one field blank to be analyzed for VOCs by EPA 624. Samples specific to JC17397 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|------------|-------------|----------------|-------------|
| JC17397-1 | Trip Blank | Groundwater | 3/31/2016 | 8:30 |
| JC17397-2 | Field | Groundwater | 3/31/2016 | 9:45 |
| JC17397-3 | MW-5 | Groundwater | 3/31/2016 | 10:00 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey the same day as collected. As indicated by the analytical laboratory sample receipt summary: sample coolers were received with temperatures below 6°C, VOAs to be analyzed for VOCs did not contain headspace, sufficient sample volume was received, custody seals were intact and containers were accurately labeled and reflected the chain of custody.

2.12.1 VOCs

Preservation and Holding Time

Samples were received within the required temperature range and were analyzed for VOCs between five and six days after collection. VOAs were preserved with hydrochloric acid (HCl) and was verified to be pH<2 by the laboratory. Samples were analyzed within method holding time.

GC/MS Instrument Performance Check

lon abundances were averaged over three scans and background was subtracted from less than 20 scans away. BFB ion abundances were within CLP guidance ranges. BFB was analyzed at the beginning of every analytical sequence and once per day (EPA 624 method).

Initial Calibration

RRF and %RSD were within control limits for the calibration used to quantitate target analytes in SDG JC17397.

Initial Calibration Verification

An ICV was analyzed after the last ICAL standard. The ICV RRF for target analytes was greater than the minimum RRF. The %D between the ICAL mean RRF and the ICV RRF was within limits of the CLP guidance for trace volatiles.

Continuing Calibration Verification

Target analytes were within %D and RT limits for opening CCVs. RRFs were greater than minimums presented in Table 4 of the CLP guidance.

Blanks

No target analytes were detected in the trip blank or method blanks. Tetrachloroethene was detected in the field blank. The concentration of tetrachloroethene in the environmental samples within SDG JC17397 was almost 1000 times greater than the concentration found in the field blank. This contamination is not expected to influence the environmental samples concentration of tetrachloroethene.

Surrogates

Surrogate recoveries in environmental samples and quality control samples were within control limits.

Laboratory Control Samples

The recovery of target analytes was within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

The matrix spike had recoveries of target analytes within control limits.

Laboratory Duplicate

A laboratory duplicate was analyzed as part of the laboratory quality control samples. The source of the laboratory duplicate was not part of SDG JC17397; all RPD values were within control limits except for chloroethane, which was detected at less than the reporting limit. This value is not controlled.

Internal Standards

The retention time of the internal standards was compared against daily beginning continuing calibration verifications. The retention times did not differ more than 10 seconds and area counts were within 50-200% of the check standard.

Target Analyte Identification

Reported target analytes are accurately identified based on ion abundances, analyte spectrums and retention times.

2.13 SDG JC22892

Data validation and review was conducted for soil and blank water samples within SDG JC22892 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 and NYDEC Analytical Services Protocol. The SDG included ten soil samples, one trip blank and one field blank to be analyzed for VOCs by EPA 8260C and percent moisture by SM2540. Trip and field blanks were sampled to show there was no contamination from transportation or field activities. Samples specific to JC22892 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|------------|---------------|--------|-------------|-------------|
| JC22892-1 | MW-9 (19.4') | Soil | 6/22/2016 | 10:45 |
| JC22892-2 | MW-9 (30′) | Soil | 6/22/2016 | 10:45 |
| JC22892-3 | MW-10 (19.5') | Soil | 6/22/2016 | 12:05 |
| JC22892-4 | MW-10 (30') | Soil | 6/22/2016 | 12:05 |
| JC22892-5 | MW-8 (19.5') | Soil | 6/22/2016 | 14:45 |
| JC22892-6 | MW-8 (30') | Soil | 6/22/2016 | 14:45 |
| JC22892-7 | MW-5I (20') | Soil | 6/23/2016 | 8:35 |
| JC22892-8 | MW-5I (23') | Soil | 6/23/2016 | 8:45 |
| JC22892-9 | MW-5I (31') | Soil | 6/23/2016 | 8:55 |
| JC22892-10 | MW-5I (50') | Soil | 6/23/2016 | 9:40 |
| JC22892-11 | FB | Soil | 6/22/2016 | 13:00 |
| JC22892-12 | Trip Blank | Soil | 6/20/2016 | 10:30 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: the cooler temperatures upon receipt were $<6^{\circ}$ C, custody seals were intact, containers were accurately labeled and reflected the chain of custody and there was no headspace in VOA (volatile organic analysis) containers. The QC blank samples were preserved with hydrochloric acid (HCI) and pH checked by the laboratory. Soil samples were collected with EnCore samplers.

2.13.1 VOCs

Holding Times

Samples were analyzed within holding times for both water and soil samples.

Instrument Performance Check

Bromofluorobenzene (BFB) was evaluated for mass spectra requirements and frequency to ensure adequate instrument performance. BFB standards were analyzed at the beginning of every sequence and within 12 hours of the final sample injected. Ion abundances were within control limits, which were averaged over three scans within the BFB peak and background subtracted from less than 20 scans away.

Initial Calibration

The calibration included at least five calibration points, was performed prior to ICVs, samples and blanks, and was within twelve hours of a BFB check standard. Relative Response Factors (RRF) and Percent Relative Standard Deviations (%RSD) were calculated for each analyte and surrogate. The mean RRF and %RSD for target analytes were within control limits.

Initial Calibration Verification

A second source calibration verification was analyzed after the soil and water calibrations. The percent difference between the mean RRF of the initial calibration and the initial calibration verification were within control limits stated by the CLP Guidance.

Continuing Calibration Verification

Continuing calibration verifications were evaluated to determine instrument performance and calibration validity during each analytical run. The percent difference between the mean RRF of the initial calibration and the CCV were within control limits stated by the CLP Guidance with the exception of the %D for carbon tetrachloride in the CCV ran 6/29/16 at 7:16 am. The affected sample is MW-5I (20'), which was ran in the sequence after the CCV. Carbon tetrachloride is considered an estimate in this sample and is UJ qualified in the data table. The %D for acetone was also outside of CLP guidance limits for the CCV ran 6/25/16 at 12:25 pm. Affected samples are MW-9 (30'), MW-8 (19.5'), MW-5I (31'), MW-5I (50') and MW-10 (30'). Acetone was not detected in these samples and results are UJ qualified as estimated values.

Blanks

Method blanks, field blanks and trip blanks were analyzed to assess possible contamination from the laboratory or the field. One method blank was analyzed per analytical batch. There were no detections in any blanks associated with

this SDG.

Surrogates

Surrogate recoveries were within control limits.

Laboratory Control Sample

One laboratory control sample was analyzed per analytical batch included in the SDG. The percent recoveries for target analytes were within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

One matrix spike was analyzed per analytical batch. Recoveries of Target analytes were within control limits.

Laboratory Duplicate

One laboratory duplicate was analyzed per analytical batch. Sample JC22892-2 (MW-9 (30')) was used as the source of the laboratory duplicate ran 6/25/16. RPD values were outside of control limits due to low analyte concentrations making it difficult to reproduce results within precision guidelines. No analytes were qualified due to the inability to accurately control the results.

Internal Standard

Required internal standards were added to samples, blank samples and quality control samples at the same specified concentration to monitor system performance and quantify target analytes. Raw data was checked against acceptance criteria for retention times and area response. Retention times did not differ more than 10 seconds from the check standard (beginning calibration check verification). Area responses were within 50-200% of the check standard.

Target Analyte Identification

Reported target analytes are accurately identified based on ion abundances, analyte spectrums and retention times.

2.13.2 Percent Moisture

Percent moisture was performed and calculated for soil samples of SDG JC22892. Samples results were verified by recalculation of raw data, where; *Percent Moisture* = $\frac{Wet Weight-Dry weight}{Wet Weight-Tare Weight} x100\%$.

2.14 SDG JC23746

Data validation and review was conducted for groundwater and blank water samples within SDG JC23746 in accordance with the EPA *National Functional Guidelines for Superfund Organic Methods Review*, January 2017 and NYDEC Analytical Services Protocol. The SDG included five groundwater samples, one trip blank and one field blank to be analyzed for VOCs by EPA 8260C. Trip and field blanks were sampled to show there was no contamination from transportation or field activities. Samples specific to JC23746 are included below.

| SDG ID | Sample ID | Matrix | Sample Date | Sample Time |
|-----------|-----------|-------------|-------------|-------------|
| JC23746-1 | MW-5I | Groundwater | 7/8/2016 | 13:19 |
| JC23746-2 | MW-8 | Groundwater | 7/8/2016 | 11:00 |
| JC23746-3 | MW-9 | Groundwater | 7/8/2016 | 12:40 |
| JC23746-4 | MW-10 | Groundwater | 7/8/2016 | 10:32 |
| JC23746-5 | MW-100 | Groundwater | 7/8/2016 | 13:19 |
| JC23746-6 | Field | Water | 7/8/2016 | 8:50 |
| JC23746-7 | Trip | Water | 7/5/2016 | 14:00 |

Samples were received by the analytical laboratory, SGS Accutest of Dayton, New Jersey. As indicated by the analytical laboratory sample receipt summary: the cooler temperatures upon receipt were $<6^{\circ}$ C, custody seals were intact, containers were accurately labeled and reflected the chain of custody and there was no headspace in VOA (volatile organic analysis) containers. The water samples were preserved with hydrochloric acid (HCI) and pH checked by the laboratory.

2.14.1 VOCs

Holding Times

Samples were analyzed between three and six days after collection; within holding time.

Instrument Performance Check

Bromofluorobenzene (BFB) was evaluated for mass spectra requirements and frequency to ensure adequate instrument performance. BFB standards were analyzed at the beginning of every sequence and within 12 hours of the final sample injected. Ion abundances were within control limits, which were averaged over three scans within the BFB peak and background subtracted from less than 20 scans away.

Initial Calibration

The calibration included at least five calibration points, was performed prior to ICVs, samples and blanks, and was within twelve hours of a BFB check standard. Relative Response Factors (RRF) and Percent Relative Standard Deviations (%RSD) were calculated for each analyte and surrogate. The mean RRF and %RSD for target analytes were within control limits.

Initial Calibration Verification

The percent difference between the mean RRF of the initial calibration and the initial calibration verification were within control limits stated by the CLP Guidance.

Continuing Calibration Verification

Continuing calibration verifications were evaluated to determine instrument performance and calibration validity during each analytical run. The percent difference between the mean RRF of the initial calibration and the CCV were within control limits stated by the CLP Guidance.

Blanks

Method blanks, field blanks and trip blanks were analyzed to assess possible contamination from the laboratory or the field. One method blank was analyzed per analytical batch. There were no detections in any blanks associated with this SDG.

Surrogates

Surrogate recoveries were within control limits.

Laboratory Control Sample

One laboratory control sample was analyzed per analytical batch included in the SDG. The percent recoveries for target analytes were within laboratory control limits.

Matrix Spike/Matrix Spike Duplicate

One MS and MSD were analyzed per analytical batch. Recoveries and RPD values for target analytes were within control limits.

Internal Standard

Required internal standards were added to samples, blank samples and quality control samples at the same specified concentration to monitor system performance and quantify target analytes. Raw data was checked against acceptance criteria for retention times and area response. Retention times did not differ more than 10 seconds from the check standard (beginning calibration check verification). Area responses were within 50-200% of the check standard.

Target Analyte Identification

Reported target analytes are accurately identified based on ion abundances, analyte spectrums and retention times.

ATTACHMENT A

RESUME OF DATA VALIDATOR



Kelsi Evans Staff Chemist

Kelsi Evans is a staff chemist with 4 years of experience in data quality and management. Beginning as an analytical chemist in an environmental laboratory, Ms. Evans can determine data quality based on standard regulatory methods. She has experience reviewing raw data produced in an environmental laboratory as a QA/QC officer and worked in hands-on positions processing and analyzing environmental samples. Her responsibilities included reviewing air, water, and soil test data produced by the laboratory for historical accuracy, appropriate use of data qualifiers, and precise identification of analytes and EPA methods. Her background in environmental analysis and data review have now cemented her position as an environmental data processor. She routinely works on sites involved in environmental sampling where screening and cleanup levels are implemented. Data packages from environmental laboratories are reviewed by Ms. Evans for criteria that determine usability. Her findings are used in reports that help ensure to clients and regulators that the data was reviewed and all possible data quality problems were discussed. Ms. Evans offers invaluable expertise in managing and interpreting environmental data involving numerous Contaminants of Concern, including metals, mercury, anions, total organic carbon, perchlorate, organochlorine pesticides, organophosphorus pesticides, PCBs, herbicides, haloacetic acids, TPH gas, TPH diesel, TPH motor oil, BTEX, and semi-volatile organics on instrumentation.

EDUCATION

B.S., Chemistry, Portland State University, Portland, Oregon, 2013

PROFESSIONAL REGISTRATIONS/ CERTIFICATION/ TRAINING

Numerous inorganic and organic chemistry instrumentation credentials Hazwoper 40-Hour Certification

GENERAL EXPERIENCE

January 2017 – Present Apex Companies, LLC, Portland, Oregon

Kelsi's role at Apex includes applying her laboratory and QA/QC experience to Apex's environmental investigation and remediation projects, including 1) laboratory data tabulation and management; 2) database development; 3) data validation and data quality review; and 4) client and regulatory specific database management.

August 2013 – December 2016 Environmental Laboratory, Northern California

QA/QC Officer and Project Manager. Ms. Evans managed a team of technicians and performed quality control review of the laboratory's air, water and soil test data for historical discrepancies to ensure appropriateness of data qualifiers before sending final reports to clients. Her responsibilities included reviewing chain of custody materials and ensuring compliance with clients' specific QAPP methods. As an analytical chemist, she worked with ion chromatography, ICP and GC instruments while managing and training other employees.

PROJECT EXPERIENCE

Port of Portland - Willamette Cove Quarterly Groundwater Monitoring and Soil Sampling, Portland, Oregon. Ms. Evans was the data manager for the December 2016 Groundwater Data Report for an Upland Facility located within the Portland Harbor Superfund Site. Her responsibilities included quality checking data tables for accuracy against laboratory reports and writing the quality assurance report. The review of the analytical data included twelve laboratory reports from multiple laboratories for VOCs, total and dissolved metals, PCB congeners, SVOCs, PAHs, anions, TOC, speciated arsenic, dioxins and furans and organochlorine pesticides in groundwater and soil. Analytical results and qualifiers were evaluated using the EPA National Functional Guidelines for Superfund Methods Data Review and individual analytical methods. Her process for reviewing the data was to read case narratives that might indicate any data quality issues, then review samples individually for holding time violations or matrix interferences (as indicated by qualifiers and surrogate recoveries) and finally quality control samples were reviewed for percent recovery and RPD control limit exceedances, blank detections and surrogate control limit failures. Sampling

documents, chain of custodies and sample receiving forms were also reviewed for sampling containers and preservation, discrepancies between bottles and chain of custody documents, unbroken chain of custody format, temperature of samples received by the laboratory and analyses. Any findings that might compromise sample data quality were documented in the Quality Assurance/Quality Control (QA/QC) review which was submitted with the Data Report to the Oregon Department of Environmental Quality.

Confidential Client - Terminal Semi-Annual Groundwater Monitoring, Portland, Oregon. Ms. Evans was the data manager for the First Semi-Annual 2017 Groundwater Monitoring Report for a bulk petroleum terminal located on the bank of the Willamette River. The environmental sampling included forty-six groundwater locations with analyses for VOCs, diesel range organics, PAHs and total and dissolved metals. Laboratory data was evaluated against the EPA National Functional Guidelines for Superfund Methods Data Review. Ms. Evans reviewed all sample handling paperwork and laboratory reports for any issues that might affect data quality. Her findings were written in a Data Quality Review report to be submitted to the Oregon Department of Environmental Quality along with the Groundwater Monitoring Report. She also reviewed and updated data qualifiers for EDD files based on valid values and data quality findings that were uploaded to a data management platform.

Gunderson LLC - Quarterly Groundwater Monitoring, Portland, Oregon. Ms. Evans managed analytical laboratory data for an industrial manufacturer located along the Willamette River. The data reviewed originated from groundwater monitoring well locations which were analyzed for VOCs, TPH diesel, PAHs, phthalates and total and dissolved metals. Laboratory data reports were evaluated for data quality based on documentation provided by the laboratory and analytical results. Any data quality issues were reviewed under the EPA National Functional Guidelines for Superfund Methods Data Review and data tables were updated accordingly to indicate any potential values that might be biased. Ms. Evans also reviewed laboratory data entry for the report data table and updated qualifiers based on her findings from the laboratory report. Exceedance ratios were calculated from screening level values and sample concentrations. These values were reviewed by Ms. Evans for accuracy based on the calculation used. Any data that might be compromised was detailed in the Analytical Laboratory Testing Quality Assurance Review that was included in the Groundwater Monitoring Report which was submitted to the Oregon Department of Environmental Quality.

Port of Portland - PDX Fire Training Facility Well Reconnaissance Sampling, Portland, Oregon. Ms. Evans facilitated communication between analytical laboratories and field staff for sampling containers and ways to reduce contamination while in the field for specialty analytes (perfluorinated compounds [PFC]). She created data tables to be included in the report, updated screening levels in tables based on site specifics and managed data to be included in the data table. Laboratory reports were reviewed by Ms. Evans for data quality based on the EPA National Functional Guidelines for Superfund Methods Data Review and all findings were included in the Quality Assurance/Quality Control (QA/QC) Review. Her review of the emerging contaminants, perfluorinated compounds, as reported were evaluated against laboratory SOPs and the EPA analytical method as basis. The EPA method was modified by the laboratory and blank detections, quality control samples and labeled standards were compared against raw data to determine usability. Field duplicates were also implemented to show field sampling and laboratory precision. During her evaluation of the analytical method and data, multiple lines of communication were opened between her and industry professionals who specialize in analytical analysis perfluorinated compounds. In addition to PFCs, groundwater samples at this location were also analyzed for VOCs, diesel and oil, PCBs, PAHs and total and dissolved metals. The review of the data was included in the Well Sampling Report.

TECHNICAL EXPERIENCE

QA/QC Officer

As a QA/QC officer at an environmental laboratory, Ms. Evans was responsible for all data quality that the laboratory produced. All raw data was evaluated by her and she determined proper use of methods, quality control completeness and overall usability. If raw data was deemed usable for a sample delivery group (SDG), she would approve the data to be reported to the client. To make these determinations she was



able to draw upon past hands-on experience and methodology provided by the EPA and Standard Methods. She maintained Environmental Laboratory Accreditation Program (ELAP) certification and submitted Performance Evaluation (PE) samples. Her data validation experience for the laboratory is as follows:

- Reviewing raw data packages and LIMS data to ensure correct final concentration based on dilution and sample preparation procedure.
- Chromatograms from GC and IC instruments were checked for ethical quantitation.
- Ensuring that sample preparation and analysis times were consistent between the LIMS and raw data submitted.
- Data submitted was compared to historical data for clients sampling locations. Any discrepancies would initiate communication between Ms. Evans and the client where it would be determined if samples needed to be prepared and analyzed again.
- Updating QC parameters in the LIMS to match client Quality Assurance Project Plans (QAPP).
- Assess data qualifiers for accuracy based on QC or sample integrity violations. Determine acceptable QC violations based on EPA methods.
- Looking at instrument sequences to determine if any carry-over was possible between samples.

Project Manager

Ms. Evans was responsible for managing projects received by the laboratory. Some of her responsibilities included:

- Reviewing sample integrity forms provided by the sample log-in department. These forms would provide a narrative of sample receiving temperature, headspace in VOAs and discrepancies between the chain of custody and bottles received.
- Ms. Evans would evaluate the amount of sample received compared to the volume needed for sample analysis and QC. She would also determine if sample containers were appropriate for methods requested.
- Sample log-in forms were checked against chain of custodies received. Ms. Evans would update any methods that were incorrectly logged in. She would communicate with clients if there were any problems with holding time, wrong matrix for method requested or incorrect preservation.

Analytical Chemist

Ms. Evans has experience working as an analytical chemist for inorganic, organic and wet chemistry departments. Her expertise in instrumental analysis helped to maintain ELAP certification by performing analysis on PE samples in all departments she worked in. Her experience as an analytical chemist includes:

- Organic analysis: GC chemist for SVOCs, pesticides (organophosphorus and organochlorine), PCBs, herbicides, TPH diesel and motor oil, TPH gas, BTEX, haloacetic acids. Duties included calibration of instrumentation. For GC/MS instruments this included tuning of instrument every 12 hours and before any run. Correct use of calibration curves (minimum number of calibration points, curve types, rho values and instrument response factors). Correctly identifying analytes whether by two-column acceptance criteria or MS ion fragments. Ethically quantitating GC chromatograms with consistency. Running calibration verification standards and instrument blanks every ten samples to ensure data quality.
- Inorganic analysis: ICP and AA chemist for metals and mercury. Calibrating each instrument at least once per day. Analyzing interference check standards to determine instrument performance. Analyzing calibration verification and instrument blanks every ten samples to ensure calibration and no carry-over.
- Other experience includes extraction technician for SVOCs, pesticides (organophosphorus and organochlorine), PCBs, herbicides, TPH diesel and motor oil, haloacetic acids, metals and mercury.

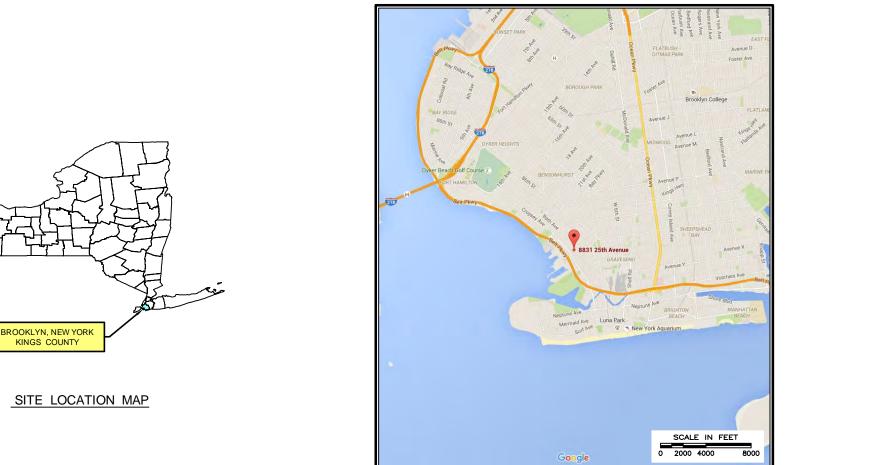


Appendix I

Garage Ventilation System Documentation

AS-BUILT DESIGN FOR GARAGE FAN SYSTEM 8831, 8841, 8851, AND 8861 BAY 25th STREET

BROOKLYN, NEW YORK





Apex Companies, LLC 223 Route 18 South, Suite 201 East Brunswick, NJ 08816 www.apexcos.com

PREPARED FOR:

8831, 8841, 8851, AND 8861 BAY 25th STREET **BROOKLYN, NEW YORK**

SHEET NO. 1 2

SCHEDULE OF DRAWINGS:

TITLE

COVER SHEET

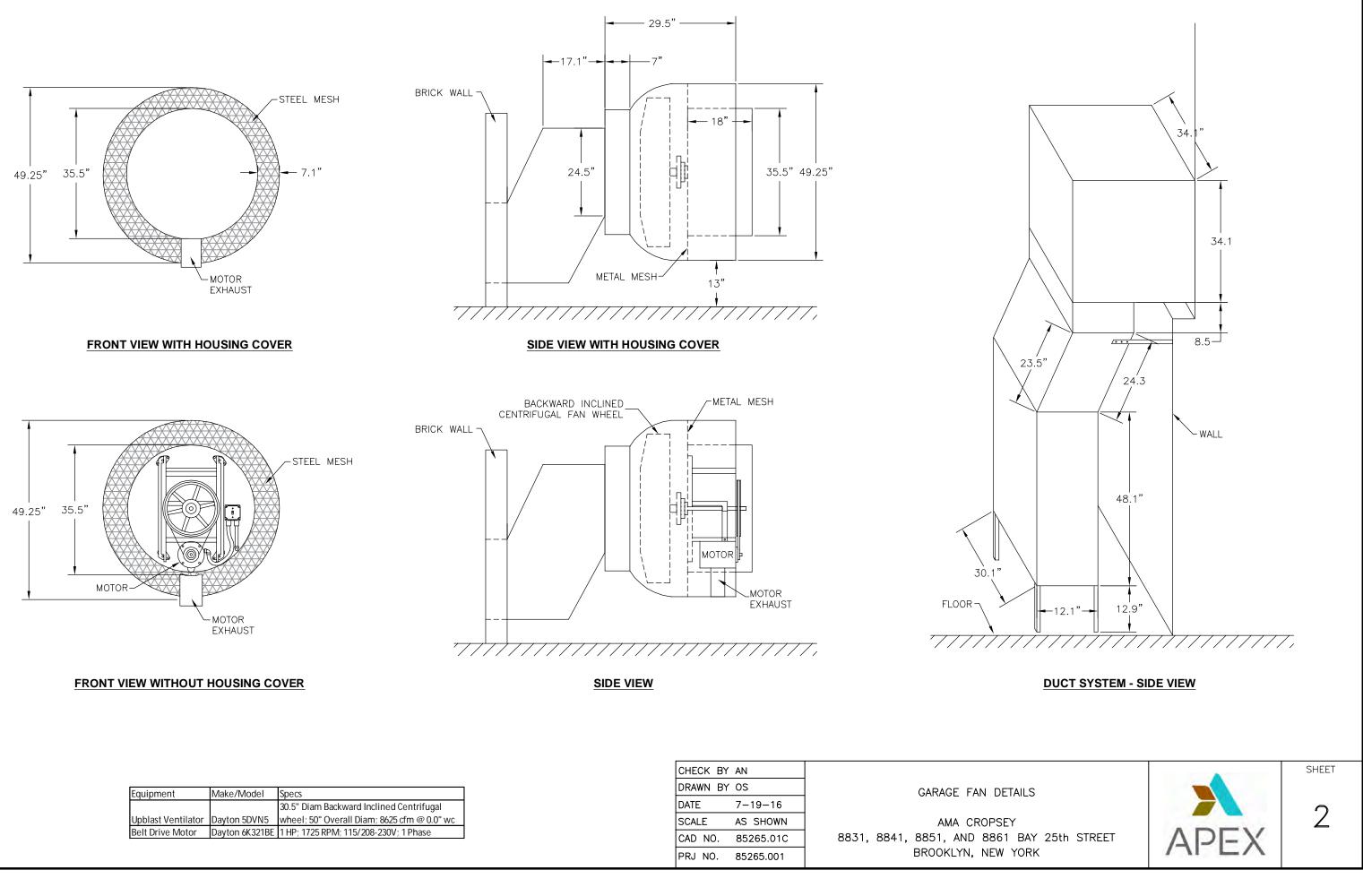
GARAGE FAN DETAILS

GARAGE FAN AIRFLOW PROCESS

GARAGE PLAN

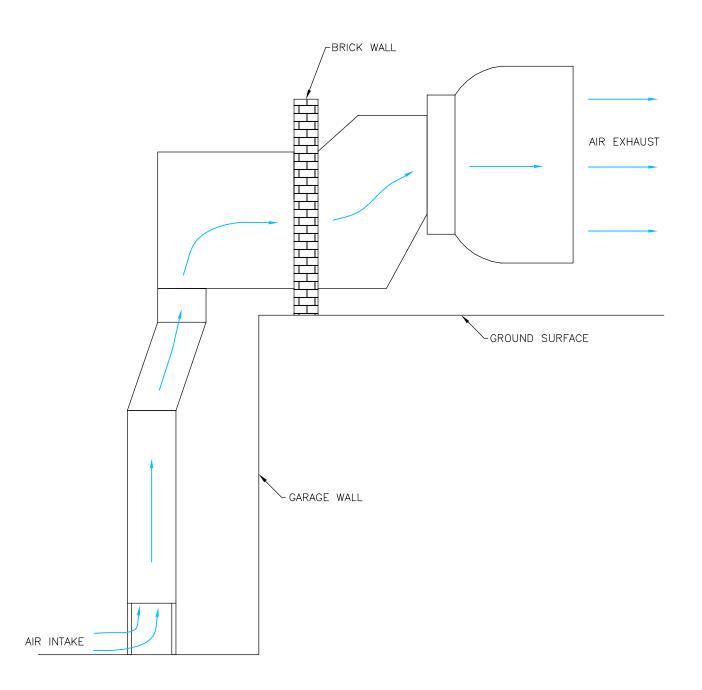


SHEET CS



| Equipment | Make/Model | Specs |
|--------------------|----------------|---|
| | | 30.5" Diam Backward Inclined Centrifugal |
| Upblast Ventilator | Dayton 5DVN5 | wheel; 50" Overall Diam; 8625 cfm @ 0.0" wc |
| Belt Drive Motor | Dayton 6K321BE | 1 HP; 1725 RPM; 115/208-230V; 1 Phase |
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| PRJ NO. | 85265.001 | BROOKLYN, NEW |
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SHEET

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