

June 13, 2018

Ms. Kerry Maloney New York State Department of Environmental Conservation 625 Broadway, 12th Floor Albany, New York 12207

RE: First Quarter Performance Monitoring Report – March 2018 432 Rodney Street
Brooklyn, New York 11211
NYSDEC BCP Site No. C224216
Langan Project No.: 170357801

Dear Ms. Maloney:

In accordance with the Site Management Plan (SMP) dated December 14, 2017, Langan Engineering, Environmental, Surveying and Landscape Architecture, D.P.C. (Langan) conducted quarterly groundwater sampling at 432 Rodney Street (the Site) located in the Borough of Brooklyn, New York. Refer to Figure 1 for the Site location. This groundwater sampling event was performed in March 2018 and represents the first quarterly sampling event for this year. This is the first quarterly sampling event since the installation and subsequent operation of a network of pressurized injection wells, and the receipt of a New York State Department of Environmental Conservation (NYSDEC) Certificate of Completion for the Site (issued December 29, 2017).

Project Background

The Site, identified as Block 2374, Lots 1, 27, 28, and 31 on the New York City Tax Map, is situated on an approximately 27,160-square-foot (±0.6235 acre) area bound by a vacant lot and residential and commercial buildings followed by Ainslie Street to the north, Keap Street to the east, Hope Street to the south, and Rodney Street to the west. The Site was most recently occupied by a packaged food storage and refrigeration facility, and has historically been used by various commercial, automotive, and light industrial companies since around 1887.

In 2016, operations ceased and the Site was remediated in accordance with the NYSDEC-approved Interim Remedial Measures Work Plan (IRMWP), dated March 22, 2016, NYSDEC-

approved IRMWP Addenda #1 and #2, dated January 9, 2017, and NYSDEC-approved Remedial Action Work Plan (RAWP), dated March 23, 2017, and approved October 31, 2017,. Under these plans, a Track 2 remedy was implemented on Lots 1 and 31, and Lots 27 and 28 follow a Track 4 remedy.

A two-phase groundwater treatment system was chosen as the remedy to treat groundwater impacted with petroleum-related volatile organic compounds (VOCs) and chlorinated volatile organic compounds (CVOCs). The treatment design included a preliminary in-situ chemical oxidation (ISCO) direct-injection of base-activated sodium persulfate, followed by the application of PlumeStop®, a liquid activated carbon substrate, through a sub-slab network of pressurized injection wells.

About 11,800 pounds of sodium hydroxide-activated sodium persulfate was direct-injected within saturated soil in the southeast corner of Lot 31 between February 21 and March 2, 2017 to treat petroleum-related VOC-impacted groundwater. Following the persulfate injection, about 56,400 pounds of PlumeStop® was applied through the well network on Lots 1 and 31 between November 8 and December 6, 2017, and about 15,600 pounds of PlumeStop® was direct-injected on Lots 27 and 28 between November 27 and December 6, 2017.

As part of the Site Management Plan (SMP) prepared by Langan and approved by the NYSDEC in December 2017, five performance monitoring wells (PMW01, PMW02, PMW03, PMW04 and PMW05) were installed on Lots 1 and 31 and two performance monitoring wells (PMW06 and PMW07) were installed on Lots 27 and 28 to monitor post-injection groundwater quality. The monitoring plan summarized in the SMP includes 1) baseline sampling, which was conducted in June 2017 (Lots 1 and 31) and November 2017 (Lots 27 and 28), after sodium hydroxide-activated sodium persulfate injections and prior to PlumeStop® injections; 2) and post-injection sampling, which is required to be conducted quarterly during the first year following the injections and semi-annually during each subsequent year as needed.

First Quarter Performance Monitoring Scope of Work

Well Purging and Sampling

Each of the seven performance monitoring wells were sampled during the first quarter 2018 event. Monitoring well sampling was conducted in accordance with the NYSDEC-approved SMP on March 20 and 23, 2018.

Each well was purged prior to sampling using the low-flow method developed by the United States Environmental Protection Agency (USEPA) ("Low-Flow [Minimal Drawdown] Ground-Water Sampling Procedures," EPA/540/S-95/504, April 1996) and accepted by the NYSDEC. Purging was performed using a peristaltic pump fitted with dedicated tubing at all wells. During purging, the turbidity, pH, temperature, conductivity, redox potential, and dissolved oxygen of



the groundwater were monitored using a Horiba U-22 Water Quality Meter with a flow-through cell. Purging was considered complete after three well volumes were purged and all parameter readings stabilized for three successive readings within a reasonable time frame. The purged water was containerized in a 55-gallon drum and temporarily stored in a secured area pending proper off-site disposal. The monitored parameters were recorded on the Well Purging and Sampling Logs provided in Attachment A.

After purging the well, a groundwater sample was collected directly from the pump discharge line using USEPA low-flow techniques at each well. For quality assurance and quality control, one field blank and a duplicate sample were collected. A trip blank was also included in each shipment for quality control. All samples were analyzed for Target Compound List (TCL) VOCs at Alpha Analytical of Westborough, Massachusetts, a New York State Department of Health (NYSDOH) Environmental Laboratory Accreditation Program (ELAP)-accredited laboratory.

Data Validation

Upon receipt of final Analytical Services Protocol (ASP) Level B laboratory reports, copies of the reports were submitted to Langan's data validation department for review in accordance with the USEPA validation guidelines for organic and inorganic data review, and the data were found to be acceptable, with no issues. There were no data flagged as either estimated or unusable.

Data reduction, validation, and reporting procedures were completed in accordance with the Quality Assurance Project Plan (QAPP) provided in Appendix I of the SMP. Data Usability Reports (DUSRs) can be found in Attachment B of this document.

First Quarter Performance Monitoring Results

Analytical Results

The laboratory analytical results for this quarterly sampling event are summarized in Table 1 and on Figure 2. Laboratory analytical reports are provided as Attachment C. Groundwater sampling results were compared to NYSDEC Technical & Operations Guidance Series (TOGS) Ambient Water Quality Standards and Guidance Values (SGVs) for Class GA water.

Two VOCs, 2-butanone and acetone, were detected above the TOGS SGVs in monitoring well PMW07 located on Lot 27. Two CVOCs, cis-1,2-dichloroethene and tetrachloroethene (PCE), were detected above the TOGS SGVs in the four monitoring wells (PM01, PM02, PM03, and PM04) located within the building footprint on Lots 1 and 31.

Based on first quarter groundwater monitoring results, the extent of the petroleum-related VOC and CVOC-impacted groundwater has decreased relative to the baseline sampling following the implementation of the two-phase groundwater treatment program. Sampling results from the baseline to the first quarter indicate a decrease of the following CVOCs:

- 1,2-dichloroethane (100%),
- cis-1,2-Dichloroethene (between 65% and 99%),



- PCE¹ (between 44% and 100%)
- trichloroethene (TCE) (between 32% and 100%, and
- vinyl chloride (between 91% and 100%);

The following petroleum-related VOC concentrations decreased, relative to the baseline concentration:

- 1,2,4-trimethylbenzene (100%),
- Benzene (100%),
- Ethylbenzene (100%),
- Isopropylbenzene (100%),
- n-propylbenzene (100%),
- p/m-xylene (100%), and
- toluene (100%).

Table 2 compares the March 2018 first quarter analytical results to the baseline sampling results collected in June and November 2017.

Closure

In general, petroleum-related VOC and CVOC concentrations detected during the June 2017 baseline sampling event have decreased by one- to two-orders of magnitude in the first quarter sampling event. Post-remediation monitoring indicates that the two-phase groundwater remedy selected for this Site appears on track to meet the remedial objective of 90% contaminant mass reduction. We recommend continued monitoring of the sub-slab performance monitoring well network on a quarterly basis, as prescribed in the SMP.

Should you have any questions, please call the undersigned at 212-479-5413.

Sincerely,

Langan Engineering, Environmental, Surveying Landscape Architecture, and Geology, D.P.C.

Michael D. Burke, PG, CHMM Principal/Vice President

¹ The PCE concentration in groundwater in PMW02 remained relatively unchanged at 15 mg/L, compared to the baseline concentration of 10 mg/L.



Enclosures:

Figure 1 Site Location Map

Figure 2 Groundwater Sampling Results Map – First Quarter March 2018

Table 1 Quarterly Groundwater Sampling Results – First Quarter 2018

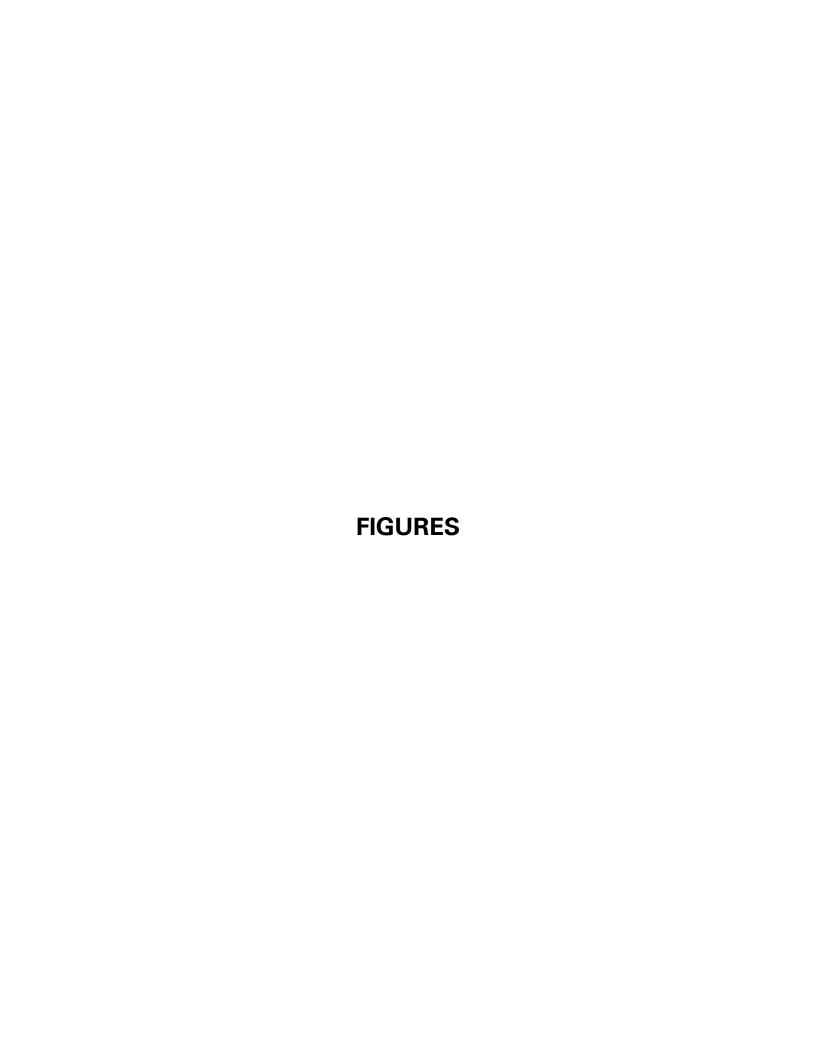
Table 2 Historical Performance Monitoring Analytical Results

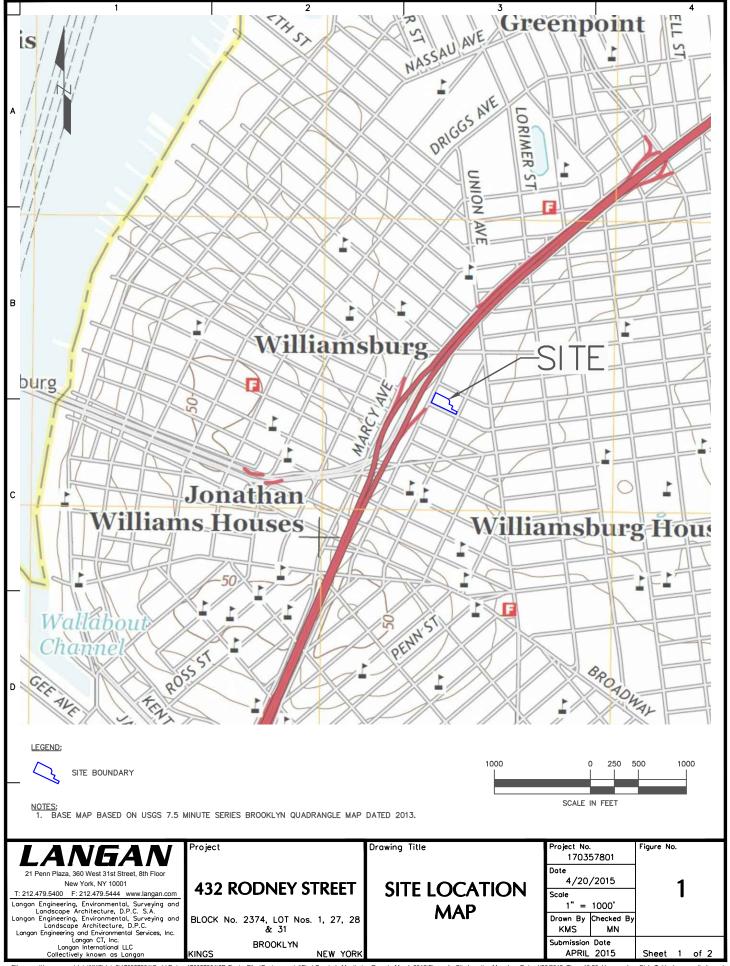
Attachment A Well Purging and Sampling Logs

Attachment B Data Usability Reports

Attachment C Laboratory Analytical Reports







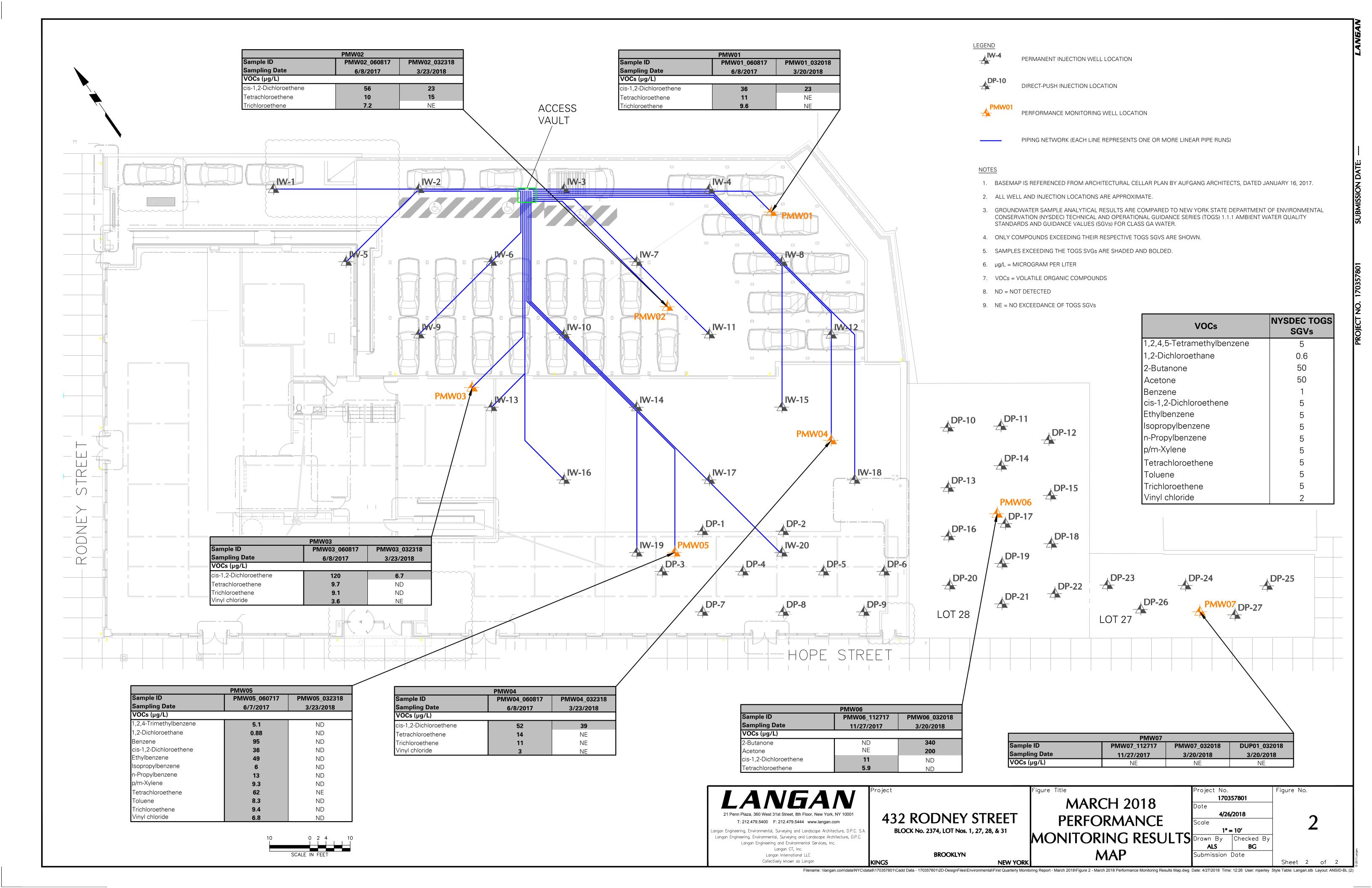




Table 1 - Quarterly Groundwater Sampling Results - First Quarter 2018 432 Rodney Street Brooklyn, New York Langan Project No. 170357801 BCP Site No. C224216

| Sample ID Sampling Date Laboratory ID | NYSDEC TOGS SGVs | PMW01_03 3/20/20 L180958 |)18 | PMW02_03 3/23/20 L1810069 | 18 | PMW03_03 3/23/20 L1810069 | 18 | PMW04_0 3/23/20 L181006 |)18 | PMW05_0 3/23/20 L181006 | 018 | PMW06_0 3/20/20 L180958 | 018 | PMW07_ 3/20/2 L18095 | 2018 | DUP01_03 3/20/20 L180958 | 018 |
|---|---------------------|--------------------------------|-----------|---------------------------------|-----------|---------------------------------|-----------|-------------------------------|-----------|-------------------------------|-----------|-------------------------------|-----------|----------------------------|-----------|--------------------------------|-----------|
| VOCs (μg/L) | | | | | | | | | | | | | | | | | |
| 1,2-Dichloroethene | ~ | 23 | | 23 | | 6.7 | | 40 | J | 2.5 | U | 2.5 | U | 2.5 | U | 2.5 | U |
| 2-Butanone | 50 | 5 | \bigcup | 5 | \bigcup | 2.3 | J | 31 | | 8.7 | | 340 | | 5 | \bigcup | 5 | \bigcup |
| Acetone | 50 | 5 | \bigcup | 5 | \bigcup | 18 | | 34 | | 22 | | 200 | | 5 | \bigcup | 5 | \bigcup |
| Benzene | 1 | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup | 0.37 | J | 0.5 | \bigcup | 0.5 | U | 0.5 | \bigcup | 0.5 | \bigcup |
| Bromodichloromethane | 50 | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup | 0.2 | J | 0.19 | J |
| Bromomethane | 5 | 2.5 | \bigcup | 2.5 | \bigcup | 1.7 | J | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup |
| Chloroform | 7 | 2.5 | \bigcup | 2.5 | \bigcup | 1.3 | J | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 3.4 | | 3.4 | |
| cis-1,2-Dichloroethene | 5 | 23 | | 23 | | 6.7 | | 39 | | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup |
| Tetrachloroethene | 5 | 3.3 | | 15 | | 0.5 | U | 2.5 | | 0.48 | J | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup |
| trans-1,2-Dichloroethene | 5 | 2.5 | \bigcup | 2.5 | U | 2.5 | \bigcup | 0.77 | J | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup |
| Trichloroethene | 5 | 2 | | 3.4 | | 0.5 | \cup | 1.6 | | 0.5 | \cup | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup |
| Vinyl chloride | 2 | 1 | \bigcup | 0.13 | J | 0.77 | J | 0.89 | J | 1 | \bigcup | 0.25 | J | 1 | \bigcup | 1 | \bigcup |

NOTES:

- 1. Groundwater sample analytical results are compared to New York State Department of Environmental Conservation (NYSDEC) Technical and Operational Guidance Series (TOGS) 1.1.1 Ambient Water Quality Standards and Guidance Values (SGVs) for Class GA water.
- 2. Concentrations exceeding NYSDEC TOGS SGVs are shaded and bold.
- 3. DUP01_032018 is a duplicate sample of PMW07_032018
- 4. μ g/L = Microgram per liter
- 5. VOCs = Volatile organic compounds
- 6. Only detected compounds are shown in the table.
- 7. \sim = Criterion does not exist.

QUALIFIERS:

- J = Detected above the Method Detection Limit (MDL) but below the Reporting Limit (RL); therefore, the result is an estimated concentration.
- U = The analyte was analyzed for, but was not detected at a level greater than or equal to the RL; the value shown in the table is the RL.

Table 2 - Historical Performance Monitoring Analytical Results

432 Rodney Street
Brooklyn, New York
Langan Project No. 170357801
BCP Site No. C224216

| Sampling Location | | | PMW01 | | | PMW02 | | | PMW03 | | | PMW04 | |
|----------------------------|------------------------------|-----------------|-------------------------------|------------------|-----------------|-------------------------------|------------------|-----------------|-------------------------------|------------------|-----------------|-------------------------------|------------------|
| Sampling Event | NYSDEC TOGS Class GA SGVs | Baseline (2016) | Post-Source Removal (2017) | Quarter 1 (2018) | Baseline (2016) | Post-Source Removal (2017) | Quarter 1 (2018) | Baseline (2016) | Post-Source Removal (2017) | Quarter 1 (2018) | Baseline (2016) | Post-Source Removal (2017) | Quarter 1 (2018) |
| Sample ID | | MW09_061616 | PMW01_060817 | PMW01_032018 | MW08S_061716 | PMW02_060817 | PMW02_032318 | MW18_070516 | PMW03_060817 | PMW03_032318 | MW10_061516 | PMW04_060817 | PMW04_032318 |
| VOCs (μg/L) | | | | | | | | 9 | • | • | | | |
| 1,1-Dichloroethane | 5 | 2.6 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| 1,1-Dichloroethene | 5 | 0.5 U | 0.2 J | 0.5 U | 0.5 U | 0.33 J | 0.5 U | <i>50</i> U | 0.56 | 0.5 U | <i>50</i> U | 0.24 J | 0.5 U |
| 1,2,4,5-Tetramethylbenzene | 5 | 2 U | 2 U | 2 U | 2 U | 2 U | 2 U | <i>200</i> U | 2 U | 2 U | <i>200</i> U | 2 U | 2 U |
| 1,2,4-Trimethylbenzene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| 1,2-Dichlorobenzene | 3 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| 1,2-Dichloroethane | 0.6 | 0.87 | 0.5 U | 0.5 U | 0.5 U | 0.14 J | 0.5 U | <i>50</i> U | 0.18 J | 0.5 U | 20 J | 0.13 J | 0.5 U |
| 1,2-Dichloroethene, Total | ~ | 99 | 36 | 23 | 66 | 56 | 23 | 4100 | 120 | 6.7 | 6500 | 52 | 40 J |
| 1,3,5-Trimethylbenzene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| 2-Butanone | 50 | 5 U | 4.7 J | 5 U | 5 U | 27 | 5 U | <i>500</i> U | 12 | 2.3 J | <i>500</i> U | 5 U | 31 |
| Acetone | 50 | 7.8 | 2.4 J | 5 U | 3 J | 28 | 5 U | <i>500</i> U | 22 | 18 | 150 J | 3.4 J | 34 |
| Benzene | 1 | 0.37 J | 0.21 J | 0.5 U | 0.6 | 0.22 J | 0.5 U | 44 J | 0.38 J | 0.5 U | 260 | 0.53 | 0.37 J |
| Bromodichloromethane | 50 | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 0.5 U | 50 U | 0.5 U | 0.5 U | 50 U | 0.5 U | 0.5 U |
| Bromomethane | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 1.4 J | 2.5 U | <i>250</i> U | 2.5 U | 1.7 J | <i>250</i> U | 2.5 U | 2.5 U |
| Chlorobenzene | 5 | 2.5 U | 1.1 J | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 1.4 J | 2.5 U |
| Chloroform | 7 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 1.3 J | <i>250</i> U | 2.5 U | 2.5 U |
| cis-1,2-Dichloroethene | 5 | 99 | 36 | 23 | 66 | 56 | 23 | 4100 | 120 | 6.7 | 6500 | 52 | 39 |
| Dichlorodifluoromethane | 5 | 5 U | 5 U | 5 U | 5 U | 5 U | 5 U | <i>500</i> U | 5 U | 5 U | <i>500</i> U | 5 U | 5 U |
| Ethylbenzene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| Isopropylbenzene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| Methyl tert butyl ether | 10 | 0.71 J | 1.1 J | 2.5 U | 1.9 J | 4.5 | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 7.8 | 2.5 U |
| Naphthalene | 10 | 0.77 J | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 0.71 J | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| n-Butylbenzene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| n-Propylbenzene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | | 2.5 U |
| o-Xylene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| p/m-Xylene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| p-Diethylbenzene | ~ | 2 U | 2 U | 2 U | 2 U | 2 U | 2 U | 200 U | 2 U | 2 U | 200 U | 2 U | 2 U |
| p-Ethyltoluene | ~ | 2 U | 2 U | 2 U | 2 U | 2 U | 2 U | 200 U | 2 U | 2 U | 200 U | 2 U | 2 U |
| sec-Butylbenzene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U |
| Tetrachloroethene | 5 | 48 | 11 | 3.3 | 27 | 10 | 15 | 500 | 9.7 | 0.5 U | 1000 | 14 | 2.5 |
| Toluene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | <i>250</i> U | | 2.5 U |
| trans-1,2-Dichloroethene | 5 | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | <i>250</i> U | 2.5 U | 2.5 U | | | 0.77 J |
| Trichloroethene | 5 | 13 | 9.6 | 2 | 5 | 7.2 | 3.4 | 130 | 9.1 | 0.5 U | 210 | 11 | 1.6 |
| Vinyl chloride | 2 | 0.5 J | 0.46 J | 1 U | 1.5 | 0.8 J | 0.13 J | 240 | 3.6 | 0.77 J | 330 | 3 | 0.89 J |
| Xylenes, Total | ~ | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 2.5 U | 250 U | 2.5 U | 2.5 U | 250 U | 2.5 U | 2.5 U |

Notes:

- 1. Groundwater sample analytical results are compared to New York State Department of Environmental Conservation (NYSDEC) Technical and Operational Guidance Series (TOGS) 1.1.1 Ambient Water Quality Standards and Guidance Values (SGVs) for Class GA drinking water.
- 2. Results exceeding the NYSDEC TOGS standards and guidance values are shaded and bolded.
- 3. Results exceeding the NYSDEC TOGS standards and guidance values but are not detected at or above the level indicated are italicized.
- 4. μg/L = micrograms per liter.
- 5. Baseline analytical results are from sampling performed in June and July 2016.
- 6. Post-source removal analytical results are from sampling performed in June and November 2017.
- 7. Quarter 1 analytical results are from sampling performed in March 2018.
- 8. VOCs = Volatile organic compounds.

Qualifiers:

- J = Analyte detected at or above the MDL (Method Detection Limit) but below the RL (Reporting Limit) data is estimated.
- U = Analyte not detected at or above the level indicated.

Table 2 - Historical Performance Monitoring Analytical Results

432 Rodney Street Brooklyn, New York Langan Project No. 170357801 BCP Site No. C224216

| Sampling Location | | | | PMW05 | j | | PMW06 PMW07 | | | | | | | | | | | |
|----------------------------|------------------------------|--------------|-----------|-------------------------|-----------|--------------|-------------|----------------|-----------|-------------------------------|--------------|-----------|---------------|-----------|------------------------------|--------|-----------|-----------|
| Sampling Event | NYSDEC TOGS Class GA SGVs | Baseline (20 | 016) | Post-Sour Removal (2 | | Quarter 1 (2 | 2018) | Baseline (2016 | 5) | Post-Source Removal (2017) | Quarter 1 (2 | 2018) | Baseline (201 | 6) | Post-Source Removal (2017 | ') | Quarter 1 | (2018) |
| Sample ID | | MW15S_061 | 1716 | PMW05_060 | 0717 | PMW05_032 | 2318 | MW12_061716 | 3 | PMW06_112717 | PMW06_03 | 2018 | MW13_06171 | 16 | PMW07_11271 | 7 | PMW07_03 | 32018 |
| VOCs (μg/L) | - | | | | | | ; | | | | | | | | | | | |
| 1,1-Dichloroethane | 5 | 100 | U | 1 | J | 2.5 | U | 12 | \bigcup | 2.5 U | 2.5 | U | 2.5 | \bigcup | 2.5 | U | 2.5 | U |
| 1,1-Dichloroethene | 5 | 20 | \bigcup | 0.5 | \bigcup | 0.5 | \bigcup | 2.5 | \cup | 0.5 U | 0.5 | \bigcup | 0.33 | J | 0.5 | \cup | 0.5 | \bigcup |
| 1,2,4,5-Tetramethylbenzene | 5 | 76 | J | 1.9 | J | 2 | \bigcup | 13 | | 2 U | 2 | \bigcup | 2 | \bigcup | 2 | \cup | 2 | \cup |
| 1,2,4-Trimethylbenzene | 5 | 2000 | | 5.1 | | 2.5 | \bigcup | 53 | | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \cup | 2.5 | \cup |
| 1,2-Dichlorobenzene | 3 | 100 | U | 0.78 | J | 2.5 | \bigcup | 12 | \bigcup | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \cup | 2.5 | \cup |
| 1,2-Dichloroethane | 0.6 | 20 | \bigcup | 0.88 | | 0.5 | \bigcup | 2.5 | \bigcup | 0.5 U | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \cup | 0.5 | \cup |
| 1,2-Dichloroethene, Total | ~ | 100 | \bigcup | 37 | J | 2.5 | \bigcup | 21 | | 11 | 2.5 | \bigcup | 46 | | 2.5 | \cup | 2.5 | \cup |
| 1,3,5-Trimethylbenzene | 5 | 580 | | 2.5 | \bigcup | 2.5 | \bigcup | 17 | | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \cup | 2.5 | \cup |
| 2-Butanone | 50 | 200 | U | 5 | \bigcup | 8.7 | | 25 | \bigcup | 5 U | 340 | | 5 | \bigcup | 5 | \cup | 5 | \cup |
| Acetone | 50 | 240 | | 42 | | 22 | | 25 | \cup | 5 U | 200 | | 4.5 | J | 5 | \cup | 5 | \cup |
| Benzene | 1 | 260 | | 95 | | 0.5 | \bigcup | 41 | | 0.69 | 0.5 | U | 0.35 | J | 0.5 | \cup | 0.5 | \bigcup |
| Bromodichloromethane | 50 | 20 | U | 0.5 | U | 0.5 | \bigcup | 2.5 | \cup | 0.5 U | 0.5 | \bigcup | 0.5 | \bigcup | 0.5 | \cup | 0.2 | J |
| Bromomethane | 5 | 100 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 12 | \cup | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | \cup | 2.5 | \cup |
| Chlorobenzene | 5 | 100 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 12 | \cup | 1 J | 2.5 | \bigcup | 3.6 | | 2.5 | \cup | 2.5 | \cup |
| Chloroform | 7 | 100 | \bigcup | 2.5 | \bigcup | 2.5 | \bigcup | 12 | U | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 3.5 | | 3.4 | |
| cis-1,2-Dichloroethene | 5 | 100 | \bigcup | 36 | | 2.5 | \bigcup | 21 | - 1 | 11 | 2.5 | \bigcup | 46 | | 2.5 | \cup | 2.5 | \cup |
| Dichlorodifluoromethane | 5 | 200 | \bigcup | 5 | U | 5 | \bigcup | 25 | U | 5 U | 5 | \bigcup | 17 | | 5 | U | 5 | \bigcup |
| Ethylbenzene | 5 | 1500 | | 49 | | 2.5 | \bigcup | 120 | | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | U | 2.5 | \bigcup |
| Isopropylbenzene | 5 | 110 | | 6 | | 2.5 | \bigcup | 20 | | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | U | 2.5 | \bigcup |
| Methyl tert butyl ether | 10 | 100 | U | 1.4 | J | 2.5 | \bigcup | 12 | \cup | 2.5 U | 2.5 | \bigcup | 1.6 | J | 2.5 | U | 2.5 | U |
| Naphthalene | 10 | 75 | J | 3.8 | | 2.5 | \bigcup | 8.5 | J | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | U | 2.5 | U |
| n-Butylbenzene | 5 | 44 | J | 2 | J | 2.5 | \bigcup | 11 | J | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | U | 2.5 | U |
| n-Propylbenzene | 5 | 410 | | 13 | | 2.5 | \bigcup | 38 | | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | U | 2.5 | U |
| o-Xylene | 5 | 2300 | | 2.3 | J | 2.5 | \bigcup | 16 | | 2.5 U | 2.5 | \bigcup | 2.5 | \bigcup | 2.5 | U | 2.5 | U |
| p/m-Xylene | 5 | 5500 | | 9.3 | | 2.5 | \bigcup | 120 | | 2.5 U | 2.5 | U | 2.5 | \bigcup | 2.5 | U | 2.5 | U |
| p-Diethylbenzene | ~ | 240 | | 0.83 | J | 2 | \bigcup | 15 | | 2 U | 2 | U | 2 | \bigcup | 2 | U | 2 | U |
| p-Ethyltoluene | ~ | 1900 | | 2 | | 2 | \bigcup | 36 | | 2 U | 2 | U | 2 | \bigcup | 2 | U | 2 | U |
| sec-Butylbenzene | 5 | 100 | U | 3 | | 2.5 | \bigcup | 9.1 | J | 0.74 J | 2.5 | U | 1.2 | J | 2.5 | U | 2.5 | U |
| Tetrachloroethene | 5 | 20 | U | 62 | | 0.48 | J | 1.8 | J | 5.9 | 0.5 | \bigcup | 67 | | 0.96 | | 0.5 | U |
| Toluene | 5 | 2300 | | 8.3 | | 2.5 | U | | J | 2.5 U | 2.5 | U | 2.5 | \bigcup | | U | 2.5 | U |
| trans-1,2-Dichloroethene | 5 | 100 | U | 0.82 | J | 2.5 | U | | U | 2.5 U | 2.5 | U | 2.5 | \bigcup | | U | 2.5 | U |
| Trichloroethene | 5 | 20 | U | 9.4 | | 0.5 | U | | | 2.9 | 0.5 | U | 8 | | 0.2 | J | 0.5 | U |
| Vinyl chloride | 2 | 40 | U | 6.8 | | 1 | U | | | 1.4 | 0.25 | J | 12 | | | U | 1 | U |
| Xylenes, Total | ~ | 7800 | | 12 | J | 2.5 | \bigcup | 140 | | 2.5 U | 2.5 | \bigcup | 2.5 | U | 2.5 | U | 2.5 | U |

Notes:

- 1. Groundwater sample analytical results are compared to New York State Department of Environmental Conservation (NYSDEC) Technical and Operational Guidance Series (TOGS) 1.1.1 Ambient Water Quality Standards and Guidance Values (SGVs) for Class GA drinking water.
- 2. Results exceeding the NYSDEC TOGS standards and guidance values are shaded and bolded.
- 3. Results exceeding the NYSDEC TOGS standards and guidance values but are not detected at or above the level indicated are italicized.
- 4. μ g/L = micrograms per liter.
- 5. Baseline analytical results are from sampling performed in June and July 2016.
- 6. Post-source removal analytical results are from sampling performed in June and November 2017.
- 7. Quarter 1 analytical results are from sampling performed in March 2018.
- 8. VOCs = Volatile organic compounds.

Qualifiers:

- J = Analyte detected at or above the MDL (Method Detection Limit) but below the RL (Reporting Limit) data is estimated.
- U = Analyte not detected at or above the level indicated.

ATTACHMENT A WELL PURGING AND SAMPLING LOGS

GROUND WATER SAMPLE FIELD INFORMATION FORM MW-1 123 Hove St Well#/Location: Job No. 170357801 03/20/18 Date: Weather: Sunny. 505 Sampling Personnel: U Kim Well Information **Purging Information** Sample ID PMW 01 - 0320 18 Low ADW Purging Method Well Depth (ft.) NIA Purging Rate (I/m; gpm) 11/14 Screened Interval (ft.) 1020 Start Purge Time NIA Casing Elevation (msl) End Purge Time 1045 2" Casing Diameter (in) Volume Purged (gal) 4.50 Depth to Water (ft.) NIA Water Elevation (msl) NIA Sampling Information Casing Volume (gal) Low Flow NIA Sampling Method 1050 PID/FID Reading (ppm) 0.0 Start Sampling Time 1055 End Sampling Time NA Depth Before Sampling (ft.) Number Bottles Collected Viale **Parameters** Dissolved Depth to рΗ Turbidity Conductivity Temp ORP **Purged Volume** Sample Time Oxygen Water (mS/cm) (NTU) (·C) (mV) (gallons) (mg/L) (ft.) 1025 7. 40 121 1.893 1.18 7:27 6.86 2.1 1.02 129 1030 D.891 7.85 10 35 6.75 0.891 0.93 130 1.7 8.05 2.75 6.68 0.892 1.7 0.83 8-22 1040 125 4.00 6.70 1045 0.892 0.62 8.05 123 4.50 Notes/Remarks Stability PH-± 0.1 unit Specific Conductance - ± 3% Temperature - ± 3% Dissolved Oxygen - ± 10% above 0.5 mg/L Turtidity - ± 10% above 5 NTU CRP/Eh-±10 millivalts Maximumflow rate - < 0.5 L/mor 0.13 gpm Maximumdrawdown-< 0.33 feet

| | GRO | UND WAT | ER SAMP | PLE FIELD II | NFORMA [*] | TION FOR | М | + |
|--|--------------|---------------|-------------|-------------------------|--|-------------------------------|------------------------|----------------------------|
| Site: 432 Koc | ney | Well#/Locatio | n: MW 2 | | Job No. | | | |
| Date: 3/23/ | 18 | | 30 5-40 | s. (lear | Sampling Per | sonnel: | . Nayo | tku |
| Well Information Sample ID Well Depth (ft) Screened interval (ft) | MAZ-C | 82318 | | Purging Informa | ation Purging Method Rate (I/m; gpm) tart Purge Time | (W.FOW - | - Deri | |
| Casing Elevation (msl) | | | | | End Purge Time | | | - |
| Casing Diameter (in) | 4 40 | | | Volu | me Purged (gal) | | | J |
| Depth to Water (ft) | 4 54 | | | | | | | |
| Water Elevation (msl) | | | | Sampling Inform | | 128 1 P176 | 11 = O9W | 1 |
| Casing Volume (gal) | 4. 5 | | 5 | | mpling Method | | | |
| PID/FID Reading (ppm) | U.U | | | | Sampling Time | 1450 | | - |
| | | | | | Sampling Time | | | - |
| | | | | | re Sampling (ft) | | | - |
| | | | | Number B | ottles Collected | 3 viais | | J |
| | | | | Dane | | | | |
| Sample Time | Temp (∘C) | рН | ORP (mV) | Conductivity (mS/cm) | Turbidity (NTU) | Dissolved Oxygen (mg/L) | Depth to Water (ft) | Purged Volume (gallons) |
| 1350 | 6.38 | 10.35 | -27 | (1.000 | 4.7 | 2.25 | - | 0.5 |
| 1355 | 6.58 | 9,21 | -14 | 0.1076 | 3.1 | 145 | | 7 |
| 1400 | (0.57 | 813 | 1 | 0.1076 | 200 | 1185 | | 1,25 |
| 1405 | 6,53 | 7.59 | -1.1 | 0.683 | 0.8 | 1.8 | | 1,75 |
| 1410 | 6.78 | 7.30 | 21 | 0,650 | 0.4 | 1.72 | | 2,125 |
| 1415 | 6.01 | 1.12 | 0.2 | 7,681 | Cil | 1:605 | | 2.325 |
| 14 80 | 7.04 | 2.02 | 42 | 0.088 | 0.0 | 1,65 | () | 7. 75 |
| 1425 | 7.12 | (0.96 | 61 | 0.690 | 0.1 | 1.64 | | 3.0 |
| 1420 | 7 05 | 6.92 | 100 | 0,690 | 0.0 | 1,70 | | 3.625 |
| 1435 | 7.12 | 6.89 | 108 | 0.686 | 010 | 1.7 | | 4.35 |
| 1490 | 7. 253 | 6.87 | 75 | 0.694 | 00 | 1.78 | | 4.625 |
| u us | 7 47 | 12,85 | 82 | 0.703 | 1.0 | 1,72 | | 5 00 |
| 1450 | 7.46 | 6,35 | 39 | 0 697 | 0.0 | 1,73 | | 5.325 |
| . 100 | | 6 / 2 / | | C.W. | 0 10 | 111/ | | |
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| | | | | | | | | |
| | | | No | tes/Remarks | | | | |
| Stability PH-± 0.1 unit Specific Conductance-± 3% Temperature-± 3% Dissolved Oxygen-±10% above 0.5 mg/L Turbicity-± 10% above 5 NTU ORP/Eh-±10 millivolts | | | 1 | | | | | |
| Maximum flow rate - < 0.5 L/ | mar 0.13 gam | | | | | | | |
| Maximum drawdown - < 0.33 | | | | | | | ~ | |

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| | GRO | JND WAT | ER SAMP | LE FIELD I | NFORMA | TION FOR | М | |
|---|-------------------------|---------------|---------|-----------------|--|-------------------|------------|---------------|
| | | | | | | | | |
| Site: | | Well#/Locatio | n: WW3 | 100 | Job No. | | | |
| Date: 3/23 | 118 | Weather: 🕚 | 05-403 | Clear | Sampling Per | sonnel: K | · · Virgo | 1/0 |
| | | | | | | | J | |
| Well Informa | rtion | | | Purging Informa | | Na.c.i. | F //3 c | 1 |
| Sample ID | | | | | Purging Method | MAN ABIA | 1- FEV) | |
| Well Depth (ft) | | | | | Rate (I/m; gpm) | 11.70 | | 4 |
| Screened interval (ft) | | | | | Start Purge Time | 1 4 4 | | - |
| Casing Elevation (msl) | | | | | End Purge Time | | <u> </u> | - |
| Casing Diameter (in) | | | | Volu | me Purged (gal) | | | _ |
| Depth to Water (ft) | | | | | enter a state of the state of t | | | |
| Water Elevation (msl) | | | | Sampling Infor | | 12/11/11 / 1 A 11 | 17. | 1 |
| Casing Volume (gal) | | | | | empling Method | 1. 1. 2 | - H/11 | - |
| PID/FID Reading (ppm) | 0.0 | | | | t Sampling Time | | 5 | - |
| | | | | | Sampling Time | | | - |
| | | | | | ore Sampling (ft) | | | 1 |
| | | | | Number B | otties Collected | 2 AMI | | 1 |
| | | | | Para | meters | | | |
| | Temp | pH | ORP | Conductivity | Turbidity | Dissolved | Depth to | Purged Volume |
| Sample Time | (∘C) | | (mV) | (mS/cm) | (NTU) | Oxygen (mg/L) | Water (ft) | (gallons) |
| 1239 | 6.39 | 11,97 | 116 | 3,43 | 5.0 | 2113 | | 1.5 |
| 1243 | 6.39 | 12,38 | 20 | 4,04 | 4.3 | 1,78 | | 1.75 |
| 1248 | 5.48 | 12.51 | -20 | 4,67 | 3,0 | 2.74 | | 2.00 |
| 1253 | 5.85 | 12.59 | -30 | 4.21 | 2.6 | 5.33 | | 2.5 |
| 1258 | 5.93 | 12.61 | -47 | 4.34 | 2.7 | 5.14 | | 2.35 |
| 1303 | 10.01 | 12.64 | -50 | 4.20 | コチ | 4.61 | | 3 |
| 1308 | 3.95 | 12.69 | - 58 | 4.35 | 3.4 | 413 | | 3.5 |
| 1313 | 5.93 | 12.77 | -67 | 4.40 | 2.5 | 2.54 | | 43.75 |
| 1315 | 5.96 | 12,73 | -73 | 4.35 | 2.7 | 3.31 | | 4 |
| 1333 | 5.78 | 12.74 | -75 | 4.40 | 2.3 | 3.22 | | 4,25 |
| 1328 | 586 | 12.77 | - 80 | 4. 25 | 218 | 3.39 | | 4.375 |
| 13.50 | 5.93 | 12.77 | -82 | 4.36 | 24 | 2,19 | | 4,75 |
| 1338 | 648 | 12,76 | 105 | 4.44 | 0.4 | 2,07 | | 5.125 |
| | | | | | | | | |
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| | | | No | tes/Remarks | L | | | L |
| Chaldille. | | | INO | ites/Remarks | | | | |
| Stablility PH-± 0.1 unit | | | | | | | | |
| Specific Conductance - ± 3% | | | | | | | | |
| | | | | | | | | |
| Temperature - ± 3% Dissolved Oxygen - ± 10% abo | an () 5 mm ⁴ | | | | | | | |
| Turbidity - ± 10% above 5 NT | | | | | | | | |
| ORP/En-±10 millivolts | • | | | | | | | |
| A.1 | mæ 0 12 ~~~ | | | | | | | |
| Azvimumflow rate -< 0.5 L/mor 0.13 gpm Azvimum drawdown -< 0.33 feet | | | | | | | | |

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| Site: | | Well#/Location | FWM :no | | Job No. | | | |
|-------------------------------|--------------|----------------|-------------|-------------------------|--------------------|-------------------------------|------------------------|--|
| Date: 3/23/2 | 018 | | - 4 4 4 | s, Clear | Sampling Pe | rsonnel: 🗸 | Nagotk | 0 |
| | | | | | | | 7 | |
| Well Informa | | | | Purging Inform | ation | · | | _ |
| Sample ID | MWH | _052318 | 1 | | Purging Method | d | | |
| Well Depth (ft) | | | 1 | Purging | Rate (I/m; gpm |) | 1. | |
| Screened Interval (ft) | | | | s | tart Purge Time | 11.70 | | |
| Casing Elevation (msl) | | | | | End Purge Time | 12:20 |) | |
| Casing Diameter (in) | | | | Volu | me Purged (gal |) | | |
| Depth to Water (ft) | | | | | | | | |
| Water Elevation (msl) | | | | Sampling Inform | mation | 1-1 | | _ |
| Casing Volume (gal) | | | | Sa | ampling Method | 1 | 14 | |
| PID/FID Reading (ppm) | 1.0 | | | Start | Sampling Time | 12:20 | | |
| | 2 | | | End | Sampling Time | | | |
| 1. | | | | Depth Befo | re Sampling (ft |) | | |
| (a |) | | | Number B | ottles Collected | d | |] |
| - | | | | | | | | |
| | | | | Para | meters | | | |
| Sample Time | Temp (∘C) | рН | ORP (mV) | Conductivity (mS/cm) | Turbidity (NTU) | Dissolved Oxygen (mg/L) | Depth to Water (ft) | Purged Volum (gallons) |
| 1120 | 7.05 | 11,85 | 50 | 0. 937 | 27.6 | 10.09 | | 0.5 |
| 1125 | 6.41 | 12.01 | 22 | 0,701 | 11,7 | 10,22 | | 1,125 |
| 1130 | 6.45 | 11,57 | 21 | 0,584 | 31.9 | 8.91 | | 250 |
| 1135 | 6.54 | 9.98 | Si | 0.590 | 41.1 | 8.04 | | 2,75 |
| 1140 | 6.72 | 8.44 | 80 | 0.687 | 19,9 | 7:37 | | 3 50 |
| 145 | 6.103 | 7,70 | 100 | 12.7110 | 15,3 | 10.94 | | 4.00 |
| 1150 | 10.39 | 7.46 | 117 | 0, 721 | 12.6 | 6.57 | | 11.25 |
| 1155 | 6.7 | 7,27 | 122 | 0.130 | 9,8 | 10 11 | | 4 625 |
| 1200 | 6 01 | 7,18 | 127 | 1) 723 | 811 | 5,810 | | 5.0 |
| 1205 | 5.99 | 7.09 | 132 | 0.744 | 1018 | 5,201 | | 675 |
| 1211) | 5.97 | 7,01 | 125 | 0.754 | 5.7 | 5.33 | | 5.5 |
| 1315 | 5.90 | 6.95 | 136 | 0.766 | 51 | 5.01 | | 5.75 |
| 1220 | 5.91 | 6.91 | 136 | | 4.2 | 4,85 | | 6.00 |
| 1 | × :- | | | C. 121 | | 7,07 | | <i>(i</i> , <i>i</i> , |
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| Stablility | | | Nc | tes/Remarks | | | | |
| H - ± 0.1 unit | | | | | | | | |
| pecific Conductance - ± 3% | 93 | | | | | | | |
| emperature - ± 3% | | | | | | | | |
| issolved Oxygen - ± 10% abov | o 0 5 m=" | | | | | | | |
| urbidity - ± 10% above 5 NTU | | | | | | | | |
| ORP/Eh - ± 10 millivolts | 8 | | | | | 2 | | |
| Maximum flow rate - < 0.5 L/m | ar 0 12 | | | | | | | |
| laximum flow rate · < 0.5 L/m | | | | | | | | |

| | GRO | UND WAT | ER SAMP | PLE FIELD II | NFORMA | TION FOR | M | |
|---|----------|----------------|--|-----------------|-------------------|------------------|--|---------------|
| Site: | | Well#/Location | - Miv/S | N. I | Job No. | | | |
| Date: 5/23/ | 18 | Weather: 3 | | SICIERY | Sampling Per | rsonnel: | K. Na. | act ko |
| | | | | / | | | N | J |
| Well Informa | | | - | Purging Informa | | - | | - |
| Sample ID | | 032318 | 1 | | Purging Method | | | |
| Well Depth (ft) | 1 | | 1 ' | | Rate (I/m; gpm) | | | |
| Screened Interval (ft) | + | | 4 / | | Start Purge Time | 1 | | _ |
| Casing Elevation (msl) | | | 4 ' | | End Purge Time | | O | |
| Casing Diameter (in) | | | 4 / | Volur | ime Purged (gal) | 1 | | |
| Depth to Water (ft) | | | 4 | | | | | |
| Water Elevation (msi) | 7 | | 4 | Sampling Inform | | - | | 7 |
| Casing Volume (gal) | 7. 1 | | 4 ! | | mpling Method | 1 1 1 1 1 | | 4 |
| PID/FID Reading (ppm) | <u> </u> | 1 | 1 1 | | t Sampling Time | | <u>) </u> | _ |
| | | | ! | | Sampling Time | | | _ |
| | | | , | | ore Sampling (ft) | | | |
| | | | , | Number Bo | ottles Collected | | > | _ |
| | | | | Para | meters | | 9) | |
| Sample Time | Temp | рН | ORP | Conductivity | Turbidity | Dissolved | Depth to | Purged Volume |
| Sample Time | (∘C) | | (mV) | (mS/cm) | (NTU) | Oxygen (mg/L) | Water (ft) | (gallons) |
| 1500 | t-22 | 7.10 | 1113 | 0.15 | 1.8 | 2.40 | | 0.75 |
| 1505 | 6.15 | 7.10 | 102 | 0.469 | 139 | 5.38 | | 1 |
| 1510 | 5.88 | 7.82 | 96 | 0.344 | 139 | 9.23 | | 1.5 |
| 1515 | 5,80 | 8,57 | 95 | 0.327 | 120 | 8,59 | | 1.625 |
| 1520 | 5,92 | 8.77 | 98 | 0.327 | 112 | 8.31 | | 1.875 |
| 1525 | 5,85 | 8,84 | 101 | 0.375 | 128 | 8.18 | | 2,75 |
| 1530 | 5,97 | 8.85 | 103 | 0,324 | 118 | 7.84 | | 2,50 |
| 15.35 | 6.14 | 8.90 | 107 | 6.321 | 117 | 7.107 | | 13.875 |
| 755AD 1540 | 6,11 | 8,91 | 104 | 0.319 | 114 | 7.45 | | 2.75 |
| TEAS MAD | | | | | | | | |
| 唐受) 1560 | | | | | | | | |
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| Stablility | | | Not | tes/Remarks | | | | |
| Statementy PH-± 0.1 unit | 0.1 | light by | 200 | and die | A | | | |
| PH-E 0.1 unit Specific Conductance-± 3% | 1 | HALL DA | NWW III | ntial disc | harose. | | | |
| Specific Conductance - ± 3% Temperature - ± 3% | 1 | N. | The state of the s | | Ú | | | |
| _ | 25 | | | | | | | |
| Dissolved Oxygen - ± 10% abo Turbidity - ± 10% above 5 NTL | - | | | | | | | |
| Turbicity - ± 10% above 5 NTC CRP/En - ± 10 millivolts | , e l | 1 | | | | | | |
| Maximumflowrate-< 0.5 L/r | | | | | | | | |
| Maximum drawdown -< 0.33 | | | | | | | | |
| | | 4 | | | | | | |

 $\textbf{Langan} \ \, \textbf{Engineering, Environmental, Surveying and Landscape Architecture, D.P.C.}$

GROUND WATER SAMPLE FIELD INFORMATION FORM 441 Keap St Site: Well#/Location: MW06 Job No. 170 357 802 Date: 03/20/18 Weather: Indoors Sampling Personnel: W. Kim **Well Information Purging Information** Sample ID PMW 06-032016 Purging Method Low Flow Well Depth (ft.) Purging Rate (I/m; gpm) NIA Screened Interval (ft.) 5-15 Start Purge Time 1320 Casing Elevation (msl) NIA End Purge Time 1405 2" Casing Diameter (in) Volume Purged (gal) Depth to Water (ft.) 10.39 Water Elevation (msl) NIA Sampling Information NIA Casing Volume (gal) Sampling Method LOW Flow PID/FID Reading (ppm) 0.0 Start Sampling Time 1410 1415 **End Sampling Time** Depth Before Sampling (ft.) NIA Number Bottles Collected 3 vials **Parameters** Dissolved Depth to рΗ Conductivity Turbidity Temp ORP Sample Time Purged Volume Oxygen Water (mS/cm) (NTU) (°C) (mV) (gallons) (mg/L) (ft.) 1325 5.50 4.17 3 281 2.88 14.03 -46 0.5 1330 59 276 4,26 3.15 13.96 -57 0.75 1335 5.77 3.55 -72 230 3.72 13.96 .00 1340 3.33 5.82 3.87 209 13.60 -76 .25 1345 7.51 6.04 138 3.15 -92 13.49 . 50 1350 6.06 2.46 142 3.03 14.84 -94 1.75 1355 6.12 3.88 121 99 2.81 14.12 3-0 1400 6.23 2.38 135 291 1410 -95 2.5 6.21 1405 2.38 2.91 14,18 3.0 128 Notes/Remarks Stability PH-± 0.1 unit PMW06_032018 (1410) Specific Conductance - ± 3% Temperature - ± 3%

Remember: Battery Connections - RED is POSITIVE and BLACK is NEGATIVE

Dissolved Oxygan - ±10% above 0.5 mg/L Turbidity - ±10% above 5 NTU CRP/En - ±10 millivolts

Maximumflow rate - < 0.5 L/mor 0.13 gpm Maximumdrawdown - < 0.33 feet

GROUND WATER SAMPLE FIELD INFORMATION FORM 441 Keap Site: St Well#/Location: MWOT Job No. 170357802 03/20/18 Indoors Date: Weather: Sampling Personnel: W. Kim **Well Information Purging Information** Sample ID PMW07_032018 Low flow Purging Method Well Depth (ft.) 15 Purging Rate (I/m; gpm) Screened Interval (ft.) 5-15 1210 Start Purge Time Casing Elevation (msl) NIA End Purge Time 1300 2 Casing Diameter (in) Volume Purged (gal) 3.75 Depth to Water (ft.) 10.45 Water Elevation (msl) NIA Sampling Information Casing Volume (gal) LOW Flow Sampling Method PID/FID Reading (ppm) 0.0 Start Sampling Time 1305 1310 End Sampling Time Depth Before Sampling (ft.) N/A Number Bottles Collected 3 rials **Parameters** Dissolved Depth to рΗ Conductivity Turbidity Temp ORP Purged Volume Sample Time Oxygen Water (mS/cm) (NTU) (·C) (mV) (gallons) (mg/L) (ft.) 1215 6.67 0.888 177 2.58 11.56 148 0.25 1220 6.53 0.857 979 1.87 13.01 145 0.50 1225 6.51 0.852 71.0 1.60 143 13.27 0.75 1230 54.2 6.49 0.852 1.41 13.45 14 5 0.80 1235 6.48 41.2 1.27 0.850 13.56 142 1.00 6.47 1240 0.846 17.9 1.06 13.74 141 1.25 1245 6.45 0.845 8.5 0.84 14.04 140 1.50 6.45 1250 0.843 5.4 0.76 140 14.09 2.25 1255 4.7 6.44 0.838 0.76 140 14.04 3.00 1300 6.44 0.839 4.1 0.77 14.09 140 3.75 Notes/Remarks Stability PMW07-032018 (1300) PH-± 0.1 unit Dupo1 -03 20 18 Specific Conductance - ± 3% collected Temperature - ± 3% Dissolved Oxygen - ± 10% above 0.5 mg/L Turbidity - ± 10% above 5 NTU ORP/Eh-±10 millivolts Maximumflow rate - < 0.5 L/mor 0.13 gpm

Maximum drawdown - < 0.33 feet

ATTACHMENT B DATA USABILITY REPORTS



2700 Kelly Road, Suite 200 Warrington, PA 18976 T: 215.491.6500 F: 215.491.6501 Mailing Address: P.O. Box 1569 Doylestown, PA 18901

To: Anna Schmiedicke, Langan Senior Staff Engineer

From: Emily Strake, Langan Senior Project Chemist/Risk Assessor

Date: April 18, 2018

Re: Data Usability Summary Report

For 432 Rodney Street Brooklyn, New York

Groundwater Samples Collected March 2018

Langan Project No.: 170357801

This memorandum presents the findings of an analytical data validation of the data generated from the analysis of groundwater samples collected March 20, 2018 by Langan Engineering and Environmental Services ("Langan") at 432 Rodney Street located in Brooklyn, New York. The samples were analyzed by Alpha Analytical Laboratories, Inc. located in Westborough, Massachusetts (NYSDOH ELAP registration # 11148) for volatile organic compounds (VOCs) using the analytical methods specified below.

VOCs by SW-846 Method 8260C

Table 1, below, summarizes the laboratory and client sample identification numbers, sample collection dates, and analytical parameters subject to review.

TABLE 1: SAMPLE SUMMARY

| SDG | Lab Sample ID | Client Sample ID | Sample Date | Analytical Parameters |
|----------|------------------|------------------|----------------|--------------------------|
| L1809583 | L1809583-01 | PMW06_032018 | 3/20/2018 | VOCs |
| L1809583 | L1809583-02 | PMW07_032018 | 3/20/2018 | VOCs |
| L1809583 | L1809583-03 | DUP01_032018 | 3/20/2018 | VOCs |
| L1809584 | L1809584-01 | PMW01_032018 | 3/20/2018 | VOCs |
| L1809584 | L1809584-02 | TB01_032018 | 3/20/2018 | VOCs |
| L1809584 | L1809584-03 | FB01_032018 | 3/20/2018 | VOCs |
| L1810069 | L1810069-01 | PMW02_032318 | 3/23/2018 | VOCs |
| L1810069 | L1810069-02 | PMW03_032318 | 3/23/2018 | VOCs |
| L1810069 | L1810069-03 | PMW04_032318 | 3/23/2018 | VOCs |
| L1810069 | L1810069-04 | PMW05_032318 | 3/23/2018 | VOCs |
| L1810069 | L1810069-05 | TB02_032318 | 3/23/2018 | VOCs |

Data Usability Summary Report For 432 Rodney Street Brooklyn, New York Langan Project No.: 170357801 April 18, 2018 Page 2 of 7

VALIDATION OVERVIEW

This data validation was performed in accordance with USEPA Region II Standard Operating Procedure (SOP) #HW-33A, "Low/Medium Volatile Data Validation" (September 2016, Revision 1), USEPA Region II SOP #HW-34A, "Trace Volatile Data Validation" (September 2016, Revision 1), the USEPA Contract Laboratory Program "National Functional Guidelines for Organic Superfund Methods Data Review" (USEPA-540-R-2017-002, January 2017) and the specifics of the methods employed.

Validation includes evaluation of the analytical data to verify that data are easily traceable and sufficiently complete to permit logical reconstruction by a qualified individual other than the originator. Items subject to review in this memorandum include holding times, sample preservation, instrument tuning, instrument calibration, laboratory blanks, laboratory control samples, system monitoring compounds, internal standard area counts, trip blanks, field blanks, field duplicates, target compound identification and quantification, chromatograms, and overall system performance.

As a result of the review process, the following qualifiers may be assigned to the data in accordance with the USEPA's guidelines and best professional judgment:

- **R** The sample results are unusable due to the quality of the data generated because certain criteria were not met. The analyte may or may not be present in the sample.
- **J** The analyte was positively identified and the associated numerical value is the approximate concentration of the analyte in the sample.
- **UJ** The analyte was not detected at a level greater than or equal to the reporting limit (RL); however, the reported RL is approximate and may be inaccurate or imprecise.
- **U** The analyte was analyzed for, but was not detected at a level greater than or equal to the level of the RL or the sample concentration for results impacted by blank contamination.
- **NJ** The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

If any validation qualifiers are assigned these qualifiers supersede any laboratory-applied qualifiers. Data that is not qualified as a result of this data validation is considered acceptable on the basis of the items subject to review. Data that is qualified as "R" are not sufficiently valid and technically supportable to be used for data interpretation. Data that is otherwise qualified due to minor data quality anomalies are usable, as qualified.



Data Usability Summary Report For 432 Rodney Street Brooklyn, New York Langan Project No.: 170357801 April 18, 2018 Page 3 of 7

TABLE 2: VALIDATOR-APPLIED QUALIFICATION

| Project Sample ID | Analysis | Analyte | CAS No. | Validator Qualifier |
|-------------------|----------|-------------------------|----------|------------------------|
| PMW06_032018 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| PMW06_032018 | SW8260C | Bromomethane | 74-83-9 | UJ |
| PMW06_032018 | SW8260C | Chloroethane | 75-00-3 | UJ |
| PMW06_032018 | SW8260C | Chloromethane | 74-87-3 | UJ |
| PMW06_032018 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| PMW06_032018 | SW8260C | Hexachlorobutadiene | 87-68-3 | UJ |
| PMW06_032018 | SW8260C | Vinyl Chloride | 75-01-4 | J |
| PMW07_032018 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| PMW07_032018 | SW8260C | Bromomethane | 74-83-9 | UJ |
| PMW07_032018 | SW8260C | Chloroethane | 75-00-3 | UJ |
| PMW07_032018 | SW8260C | Chloromethane | 74-87-3 | UJ |
| PMW07_032018 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| PMW07_032018 | SW8260C | Hexachlorobutadiene | 87-68-3 | UJ |
| PMW07_032018 | SW8260C | Vinyl Chloride | 75-01-4 | UJ |
| DUP01_032018 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| DUP01_032018 | SW8260C | Bromomethane | 74-83-9 | UJ |
| DUP01_032018 | SW8260C | Chloroethane | 75-00-3 | UJ |
| DUP01_032018 | SW8260C | Chloromethane | 74-87-3 | UJ |
| DUP01_032018 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| DUP01_032018 | SW8260C | Hexachlorobutadiene | 87-68-3 | UJ |
| DUP01_032018 | SW8260C | Vinyl Chloride | 75-01-4 | UJ |
| PMW01_032018 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| PMW01_032018 | SW8260C | Bromomethane | 74-83-9 | UJ |
| PMW01_032018 | SW8260C | Chloroethane | 75-00-3 | UJ |
| PMW01_032018 | SW8260C | Chloromethane | 74-87-3 | UJ |
| PMW01_032018 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| PMW01_032018 | SW8260C | Hexachlorobutadiene | 87-68-3 | UJ |
| PMW01_032018 | SW8260C | Vinyl Chloride | 75-01-4 | UJ |
| TB01_032018 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| TB01_032018 | SW8260C | Bromomethane | 74-83-9 | UJ |
| TB01_032018 | SW8260C | Chloroethane | 75-00-3 | UJ |



Data Usability Summary Report For 432 Rodney Street Brooklyn, New York Langan Project No.: 170357801 April 18, 2018 Page 4 of 7

| Project Sample ID | Analysis | Analyte | CAS No. | Validator Qualifier |
|-------------------|----------|---------------------------|------------|------------------------|
| TB01_032018 | SW8260C | Chloromethane | 74-87-3 | UJ |
| TB01_032018 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| TB01_032018 | SW8260C | Hexachlorobutadiene | 87-68-3 | UJ |
| TB01_032018 | SW8260C | Vinyl Chloride | 75-01-4 | UJ |
| FB01_032018 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| FB01_032018 | SW8260C | Bromomethane | 74-83-9 | UJ |
| FB01_032018 | SW8260C | Chloroethane | 75-00-3 | UJ |
| FB01_032018 | SW8260C | Chloromethane | 74-87-3 | UJ |
| FB01_032018 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| FB01_032018 | SW8260C | Hexachlorobutadiene | 87-68-3 | UJ |
| FB01_032018 | SW8260C | Vinyl Chloride | 75-01-4 | UJ |
| PMW02_032318 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| PMW02_032318 | SW8260C | Bromochloromethane | 74-97-5 | UJ |
| PMW02_032318 | SW8260C | Bromomethane | 74-83-9 | UJ |
| PMW02_032318 | SW8260C | Chloroform | 67-66-3 | UJ |
| PMW02_032318 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| PMW02_032318 | SW8260C | Trichloroethylene (TCE) | 79-01-6 | J |
| PMW02_032318 | SW8260C | Trans-1,3-Dichloropropene | 10061-02-6 | UJ |
| PMW03_032318 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| PMW03_032318 | SW8260C | Bromochloromethane | 74-97-5 | UJ |
| PMW03_032318 | SW8260C | Bromomethane | 74-83-9 | J |
| PMW03_032318 | SW8260C | Chloroform | 67-66-3 | J |
| PMW03_032318 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| PMW03_032318 | SW8260C | Trichloroethylene (TCE) | 79-01-6 | UJ |
| PMW03_032318 | SW8260C | Trans-1,3-Dichloropropene | 10061-02-6 | UJ |
| PMW04_032318 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| PMW04_032318 | SW8260C | Bromochloromethane | 74-97-5 | UJ |
| PMW04_032318 | SW8260C | Bromomethane | 74-83-9 | UJ |
| PMW04_032318 | SW8260C | Chloroform | 67-66-3 | UJ |
| PMW04_032318 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| PMW04_032318 | SW8260C | Trichloroethylene (TCE) | 79-01-6 | J |
| PMW04_032318 | SW8260C | Trans-1,3-Dichloropropene | 10061-02-6 | UJ |



Data Usability Summary Report For 432 Rodney Street Brooklyn, New York Langan Project No.: 170357801 April 18, 2018 Page 5 of 7

| Project Sample ID | Analysis | Analyte | CAS No. | Validator Qualifier |
|-------------------|----------|---------------------------|------------|------------------------|
| PMW05_032318 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| PMW05_032318 | SW8260C | Bromochloromethane | 74-97-5 | UJ |
| PMW05_032318 | SW8260C | Bromomethane | 74-83-9 | UJ |
| PMW05_032318 | SW8260C | Chloroform | 67-66-3 | UJ |
| PMW05_032318 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| PMW05_032318 | SW8260C | Trichloroethylene (TCE) | 79-01-6 | UJ |
| PMW05_032318 | SW8260C | Trans-1,3-Dichloropropene | 10061-02-6 | UJ |
| TB02_032318 | SW8260C | 1,4-Dioxane (P-Dioxane) | 123-91-1 | UJ |
| TB02_032318 | SW8260C | Bromochloromethane | 74-97-5 | UJ |
| TB02_032318 | SW8260C | Bromomethane | 74-83-9 | UJ |
| TB02_032318 | SW8260C | Chloroform | 67-66-3 | UJ |
| TB02_032318 | SW8260C | Dichlorodifluoromethane | 75-71-8 | UJ |
| TB02_032318 | SW8260C | Trichloroethylene (TCE) | 79-01-6 | UJ |
| TB02_032318 | SW8260C | Trans-1,3-Dichloropropene | 10061-02-6 | UJ |

MAJOR DEFICIENCIES:

Major deficiencies include those that grossly impact data quality and necessitate the rejection of results. No major deficiencies were identified.

MINOR DEFICIENCIES:

Minor deficiencies include anomalies that directly impact data quality and necessitate qualification, but do not result in unusable data. The section below describes the minor deficiencies that were identified.

VOCs by USEPA Method 8260C:

The initial calibration (ICAL) analyzed for instrument VOA101 exhibited a low average response factor (RF) for 1,4-dioxane (0.001). The associated results in samples DUP01_032018, PMW06_032018, PMW07_032018, TB01_032018, FB01_032018 and PMW01_032018 are qualified as "UJ".

The initial calibration verification (ICV) analyzed on 3/16/2018 at 1:14 a.m. exhibited percent differences (%Ds) greater than the control limit for dichlorodifluoromethane (69.2%), chloromethane (49.3%), vinyl chloride (44.9%), bromomethane (39.8%), chloroethane (31.3%)



Data Usability Summary Report For 432 Rodney Street Brooklyn, New York

Langan Project No.: 170357801 April 18, 2018 Page 6 of 7

and 1,4-dioxane (22.7%). The associated results in samples DUP01_032018, PMW06_032018,

PMW07_032018, TB01_032018, FB01_032018 and PMW01_032018 are qualified as "J" or

"UJ" based on potential indeterminate bias.

The continuing calibration verification (CCV) analyzed on 3/25/2018 at 9:36 exhibited %Ds

greater than the control limit for 1,4-dioxane (39.4%) and hexachlorobutadiene (22.4%). The

associated results in samples DUP01_032018, PMW06_032018, PMW07_032018,

TB01_032018, FB01_032018 and PMW01_032018 are qualified as "UJ" based on potential

indeterminate bias.

The laboratory control sample/laboratory control sample duplicate (LCS/LCSD) for batch

WG1101340 exhibited a relative percent difference (RPD) greater than the control limit for 1,4-

dioxane (24%); the associated results in samples PMW02_032318, PMW03_032318.

PMW04_03218, PMW05_032318 and TB02_032318 are qualified as "UJ" based on potential

indeterminate bias.

The ICAL analyzed for instrument GONZO exhibited a low average RF for 1,4-dioxane (0.001);

the associated results in samples PMW02_032318, PMW03_032318. PMW04_03218,

PMW05_032318 and TB02_032318 are qualified as "UJ".

The ICV analyzed on 2/15/2018 at 5:16 exhibited %Ds greater than the control limit for

dichlorodifluoromethane (45.1%), bromochloromethane (-20.2%), chloroform (-20.5%),

trichloroethene (-21.9%) and trans-1,3-dichloropropene (-20.9%); the associated results in

samples PMW02_032318, PMW03_032318. PMW04_03218, PMW05_032318 and

TB02_032318 are qualified as "UJ" based on potential indeterminate bias.

The CCV analyzed on 3/28/2018 at 8:34 exhibited %Ds greater than the control limit for

bromomethane (-49.1%) and bromochloromethane (-22.8%); the associated results in samples

PMW02_032318, PMW03_032318. PMW04_03218, PMW05_032318 and TB02_032318 are

qualified as "J" or "UJ" based on potential indeterminate bias.

OTHER DEFICIENCIES:

Other deficiencies include anomalies that do not directly impact data quality and do not

necessitate qualification. The section below describes the other deficiencies that were

identified.

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Data Usability Summary Report For 432 Rodney Street Brooklyn, New York Langan Project No.: 170357801

April 18, 2018 Page 7 of 7

VOCs by USEPA Method 8260C:

The CCV analyzed on 3/26/2018 at 8:36 exhibited %Ds greater than the control limit for chloromethane (-36.3%), 1,1-dichloroethene (-21.6%), and 1,4-dioxane (37.9%). The associated

results were reported from the initial analysis of the sample; no qualification is necessary.

COMMENTS:

One field duplicate and parent sample pair (PMW07_032018 & DUP01_032018) was collected

and analyzed for all parameters. For results less than 5X the RL, analytes meet the precision

criteria if the absolute difference is less than ±1X the RL. For results greater than 5X the RL,

analytes meet the precision criteria if the RPD is less than or equal to 30%. All parameters met

the precision criteria.

On the basis of this evaluation, the laboratory appears to have followed the specified analytical

methods with the exception of errors discussed above. If a given fraction is not mentioned

above, that means that all specified criteria were met for that parameter. All laboratory data

packages met ASP Category B requirements and all sample holding times were met.

All data are considered usable, as qualified. In addition, completeness, defined as the

percentage of analytical results that are judged to be valid, is 100%.

Signed:

Emily Strake, CEP

Senior Project Chemist/Risk Assessor

ATTACHMENT C LABORATORY ANALYTICAL REPORTS



ANALYTICAL REPORT

Lab Number: L1809583

Client: Langan Engineering & Environmental

21 Penn Plaza

360 W. 31st Street, 8th Floor New York, NY 10001-2727

ATTN: Brian Gochenaur Phone: (212) 479-5590

Project Name: 441 KEAP STREET

Project Number: 170357802

Report Date: 03/28/18

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), NJ NELAP (MA935), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-14-00197).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



03/20/18

Project Name: 441 KEAP STREET

DUP01_032018

Project Number: 170357802

L1809583-03

 Lab Number:
 L1809583

 Report Date:
 03/28/18

03/20/18 00:00

Collection Alpha Sample Sample ID Date/Time **Receive Date** Location Client ID Matrix 03/20/18 14:10 WATER BROOKLYN, NEW YORK 03/20/18 L1809583-01 PMW06_032018 WATER BROOKLYN, NEW YORK 03/20/18 13:00 03/20/18 L1809583-02 PMW07_032018

BROOKLYN, NEW YORK

WATER



L1809583

Lab Number:

Project Name: 441 KEAP STREET

Project Number: 170357802 **Report Date:** 03/28/18

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. All specific QC information is also incorporated in the Data Usability format of our Data Merger tool where it can be reviewed along with any associated usability implications. Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances the specific failure is not narrated but noted in the associated QC table. The information is also incorporated in the Data Usability format of our Data Merger tool where it can be reviewed along with any associated usability implications.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

HOLD POLICY

For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Client Service Representative and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

| Please | contact | Client | Services | at 800- | -624-9220 | with a | nv c | nuestions |
|--------|---------|--------|-----------|---------|-----------|--------|------|-------------------------|
| loase | Contact | Olicit | OCI VICCO | at ooo | 02- 0220 | with a | ıy c | _f ucsiloris. |



Serial_No:03281812:33

L1809583

Lab Number:

Project Name: 441 KEAP STREET

Project Number: 170357802 **Report Date:** 03/28/18

Case Narrative (continued)

Report Revision

March 28, 2018: This report includes the results of the sample "DUP01_032018".

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:

Title: Technical Director/Representative Date: 03/28/18

Custen Walker Cristin Walker

ORGANICS



VOLATILES



Serial_No:03281812:33

Project Name: 441 KEAP STREET

Project Number: 170357802

SAMPLE RESULTS

Lab Number: L1809583

Report Date: 03/28/18

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Lab ID: L1809583-01 Client ID: PMW06_032018

Sample Location: BROOKLYN, NEW YORK

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 03/25/18 16:08

Analyst: AD

| Date Collected: | 03/20/18 14:10 |
|-----------------|----------------|
| Date Received: | 03/20/18 |
| Field Pren: | Not Specified |

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---------------------------------|---------------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - We | stborough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | 0.25 | J | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: 441 KEAP STREET

Project Number: 170357802

SAMPLE RESULTS

Lab Number: L1809583

Report Date: 03/28/18

Lab ID: L1809583-01

PMW06_032018 Client ID:

Sample Location: BROOKLYN, NEW YORK Date Collected: 03/20/18 14:10

Date Received: 03/20/18

Field Prep: Not Specified

Sample Depth:

| Volatile Organics by GC/MS - Westborough | h Lab | | | | | |
|--|-------|---|--------------|------|------|-------|
| Trichloroothono | | | | | | |
| | ND | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | 200 | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | 330 | E | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: Lab Number: 441 KEAP STREET L1809583

Project Number: Report Date: 170357802 03/28/18

SAMPLE RESULTS

Lab ID: L1809583-01 Date Collected: 03/20/18 14:10

Client ID: Date Received: 03/20/18 PMW06_032018

Sample Location: Field Prep: Not Specified BROOKLYN, NEW YORK

Sample Depth:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|------------------------------------|------------|-----------|-------|-----|------|-----------------|--|
| Volatile Organics by GC/MS - Westh | orough Lab | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 102 | 70-130 | |
| Toluene-d8 | 105 | 70-130 | |
| 4-Bromofluorobenzene | 106 | 70-130 | |
| Dibromofluoromethane | 98 | 70-130 | |



03/20/18 14:10

Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

SAMPLE RESULTS

Lab ID: L1809583-01 D

Client ID: PMW06_032018

Sample Location: BROOKLYN, NEW YORK Date Received: 03/20/18 Field Prep: Not Specified

Date Collected:

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/26/18 10:56

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|--|--------|-----------|-------|----|-----|-----------------|--|
| Volatile Organics by GC/MS - Westborou | gh Lab | | | | | | |
| 2-Butanone | 340 | | ug/l | 25 | 9.7 | 5 | |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 105 | | 70-130 |
| Toluene-d8 | 104 | | 70-130 |
| 4-Bromofluorobenzene | 103 | | 70-130 |
| Dibromofluoromethane | 98 | | 70-130 |



Project Name: 441 KEAP STREET

Project Number: 170357802

SAMPLE RESULTS

Lab Number: L1809583

Report Date: 03/28/18

Lab ID: L1809583-02

Client ID: PMW07_032018

Sample Location: BROOKLYN, NEW YORK

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/25/18 16:37

Analyst: ΑD

| Date Collected: | 03/20/18 13:00 |
|-----------------|----------------|
| Date Received: | 03/20/18 |
| Field Prep: | Not Specified |

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---------------------------------|---------------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - We | stborough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | 3.4 | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | 0.20 | J | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |

Project Name: 441 KEAP STREET

Project Number: 170357802

SAMPLE RESULTS

Report Date: 03/28/18

Lab ID: L1809583-02

PMW07_032018 Client ID: Sample Location: BROOKLYN, NEW YORK Date Received: Field Prep:

03/20/18 13:00 03/20/18

L1809583

Date Collected:

Lab Number:

Not Specified

Sample Depth:

| Volatile Organics by GC/MS - Westborough Lab Viol ughl 0.50 0.18 1 1.2-Olchiorobexone ND ughl 2.5 0.70 1 1.4-Olchiorobexone ND ughl 2.5 0.70 1 1.4-Olchiorobexone ND ughl 2.5 0.70 1 Methyl feet Luyl ether ND ughl 2.5 0.70 1 PmXylene ND ughl 2.5 0.70 1 Vylene, Total ND ughl 2.5 0.70 1 Vylene, Total ND ughl 2.5 0.70 1 Vylene, Total ND ughl 2.5 0.70 1 Jest-(2-Olchoroethene ND ughl 2.5 0.70 1 Jest-(2-Olchoroethene, Total ND ughl 2.5 0.70 1 Dibroordeflere ND ughl 2.5 0.70 1 Als-(2-Olchoroethene ND ughl 2.5 < | Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | | |
|--|--|--------|-----------|-------|------|------|-----------------|--|--|
| 1.2 Dichlorobenzene | Volatile Organics by GC/MS - Westborough Lab | | | | | | | | |
| 1,2-Dichlorobenzene ND ugil 2,5 0,70 1 1,3-Dichlorobenzene ND ugil 2,5 0,70 1 1,3-Dichlorobenzene ND ugil 2,5 0,70 1 Methyl terb tuyl ether ND ugil 2,5 0,70 1 o-Xylene ND ugil 2,5 0,70 1 o-Xylene ND ugil 2,5 0,70 1 dis-1,2-Dichloroethene ND ugil 2,5 0,70 1 1,2-Dichloroethene, Total ND ugil 2,5 0,70 1 Dibromomethane ND ugil 2,5 0,70 1 1,2-Dichloroethene, Total ND ugil 2,5 0,70 1 Dibromomethane ND ugil 2,5 0,70 1 Actychirline ND ugil 2,5 0,70 1 Syrene ND ugil 2,5 0,70 1 <td>Trichloroethene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>0.50</td> <td>0.18</td> <td>1</td> | Trichloroethene | ND | | ug/l | 0.50 | 0.18 | 1 | | |
| 1,3-Dichlorobenzene ND ugl 2,5 0,70 1 1,4-Dichlorobenzene ND ugl 2,5 0,70 1 Methyl tert buryl ether ND ugl 2,5 0,70 1 o-Sylene ND ugl 2,5 0,70 1 o-Sylene ND ugl 2,5 0,70 1 xylenes, Total ND ugl 2,5 0,70 1 1,2-Dichloroethene ND ugl 2,5 0,70 1 1,2-Dichloroethene, Total ND ugl 2,5 0,70 1 1,2-Dichloroethene, Total ND ugl 2,0 1,0 1 1,2-Dichloroethene, Total ND ugl 2,0 1,0 1 1,2-Dichloroethene, Total ND ugl 2,0 1,0 1 2,2-Dichloroethene, Total ND ugl 2,0 1,0 1 Styrene ND ugl 2,0 1,0 1 <td>1,2-Dichlorobenzene</td> <td>ND</td> <td></td> <td>_</td> <td>2.5</td> <td>0.70</td> <td>1</td> | 1,2-Dichlorobenzene | ND | | _ | 2.5 | 0.70 | 1 | | |
| Methyl tert budyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND ug/l 2.5 0.70 1 o-Xylene ND ug/l 2.5 0.70 1 o-Xylenes ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 1,2-Dichloroethene, Total ND ug/l 2.5 0.70 1 Dibromemsthane ND ug/l 5.0 1.0 1 Acrylonkride ND ug/l 5.0 1.0 1 Acrylonkride ND ug/l 5.0 1.0 1 Styrene ND ug/l 5.0 1.0 1 Styrene ND ug/l 5.0 1.0 1 Obchtorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.0 1 Vilyi acetate | 1,3-Dichlorobenzene | ND | | | 2.5 | 0.70 | 1 | | |
| ND | 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| o-Xylene ND ug1 2.5 0.70 1 Xylenes, Total ND ug1 2.5 0.70 1 cis-1,2-Dichloroethene, Total ND ug1 2.5 0.70 1 Dibromomethane ND ug1 2.5 0.70 1 Dibromomethane ND ug1 2.5 0.70 1 Acrylontrile ND ug1 2.5 0.70 1 Acrylontrile ND ug1 2.5 0.70 1 Styrene ND ug1 5.0 1.5 1 Acetone ND ug1 5.0 1.5 1 Acetone ND ug1 5.0 1.0 1 Carbon disulfide ND< | Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Xylenes, Total ND ug/l 2.5 0.70 1 cis-1,2-Dichlorcethene ND ug/l 2.5 0.70 1 cis-1,2-Dichlorcethene, Total ND ug/l 2.5 0.70 1 Dichloromethane ND ug/l 2.5 0.70 1 L;2-Dichloroptopane ND ug/l 2.5 0.70 1 Acytonitrile ND ug/l 5.0 1.5 1 Syrene ND ug/l 5.0 1.5 1 Dichlorodfluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Vinyl acetate ND ug/l 5.0 1.0 1 | p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 1,2-Dichloroethene, Total ND ug/l 2.5 0.70 1 Dibromomethane ND ug/l 5.0 1.0 1 1,2-Trichloropropane ND ug/l 5.0 0.70 1 Acrylontrile ND ug/l 5.0 0.70 1 Styrene ND ug/l 5.0 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Styria acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 4-Hexthyl-2-pentanone ND ug/l 2.5 0.70 1 | o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,2-Dichloroethene, Total ND ug/l 2,5 0,70 1 | Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Dibromomethane ND ug/l 5.0 1.0 1 1.2.3-Trichloropropane ND ug/l 2.5 0.70 1 Acrylonitrile ND ug/l 5.0 1.5 1 Styrene ND ug/l 5.0 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.5 1 Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.0 1 Viryl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 1,3-Dichropropane ND ug/l 2.5 0.70 1 1,1,1,2-T | cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,2,3-Trichioropropane ND ug/l 2,5 0,70 1 | 1,2-Dichloroethene, Total | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Acrylonitrile ND ug/l 5.0 1.5 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.0 1 1-ynyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 1,2-Dibromochlane ND ug/l 2.5 0.70 1 1,1-1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 1,1 | Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 | | |
| Syrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.0 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 8-Methyl-2-pentanone ND ug/l 5.0 1.0 1 8-Pothachoromethane ND ug/l 2.5 0.70 1 1,2-Distromethane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 | 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Dichlorodiffluoromethane ND ug/l 5.0 1.0 1 1 1 1 1 1 1 1 1 | Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 | | |
| Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 -Butylbenzene ND ug/l 2.5 0.70 1 < | Styrene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 <t< td=""><td>Dichlorodifluoromethane</td><td>ND</td><td></td><td>ug/l</td><td>5.0</td><td>1.0</td><td>1</td></t<> | Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 | | |
| 2-Butanone ND ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 | Acetone | ND | | ug/l | 5.0 | 1.5 | 1 | | |
| Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 <td>Carbon disulfide</td> <td>ND</td> <td></td> <td>ug/l</td> <td>5.0</td> <td>1.0</td> <td>1</td> | Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 | | |
| 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 c-Chlorotoluene ND ug/l 2.5 0.70 1 | 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 | | |
| 2-Hexanone ND ug/l 5.0 1.0 1 | Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 | | |
| Bromochloromethane ND | 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 | | |
| 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.0 0.65 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 | 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 | | |
| 1,2-Dibromoethane ND ug/l 2.0 0.65 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropyltoluene ND ug/l 2.5 0.70 1 | Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 c-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropyltenue ND ug/l 2.5 0.70 1 | 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 | | |
| Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 sopropylbenzene ND ug/l 2.5 0.70 1 ug/l 2.5 0.70 1 ug/l 2.5 0.70 1 | n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| p-Isopropyltoluene ND ug/l 2.5 0.70 1 | Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| | Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Naphthalene ND ug/l 2.5 0.70 1 | p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| | Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 | | |



Project Name: Lab Number: 441 KEAP STREET L1809583

Project Number: Report Date: 170357802 03/28/18

SAMPLE RESULTS

Lab ID: L1809583-02 Date Collected: 03/20/18 13:00

Client ID: Date Received: 03/20/18 PMW07_032018

Sample Location: Field Prep: BROOKLYN, NEW YORK Not Specified

Sample Depth:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------|-----------|-------|-----|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | jh Lab | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 102 | 70-130 | |
| Toluene-d8 | 105 | 70-130 | |
| 4-Bromofluorobenzene | 104 | 70-130 | |
| Dibromofluoromethane | 97 | 70-130 | |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

SAMPLE RESULTS

Lab ID: L1809583-03 Date Collected: 03/20/18 00:00

Client ID: Date Received: 03/20/18 DUP01_032018 Sample Location: Field Prep: BROOKLYN, NEW YORK Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/25/18 15:40

Analyst: ΑD

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|----------------------------------|---------------|-----------|-------|------|------|-----------------|--|
| Volatile Organics by GC/MS - Wes | stborough Lab | | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Chloroform | 3.4 | | ug/l | 2.5 | 0.70 | 1 | |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 | |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 | |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 | |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 | |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 | 1 | |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 | |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Bromodichloromethane | 0.19 | J | ug/l | 0.50 | 0.19 | 1 | |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 | |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 | |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 | |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 | |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 | |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 | |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 | |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 | 1 | |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 | |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| | | | | | | | |



Project Name: 441 KEAP STREET

Project Number: 170357802

SAMPLE RESULTS

Lab Number: L1809583

Report Date: 03/28/18

Lab ID: L1809583-03

Client ID: DUP01_032018

Sample Location: BROOKLYN, NEW YORK Date Collected: Date Received:

Field Prep:

03/20/18 00:00 03/20/18

Not Specified

Sample Depth:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--------------------------------|-----------------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - V | Vestborough Lab | | | | | |
| Trichloroothono | ND | | // | 0.50 | 0.10 | 4 |
| Trichloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | <u> </u> |
| cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |
| | | | | | | |



Project Name: 441 KEAP STREET Lab Number: L1809583

Project Number: 170357802 **Report Date:** 03/28/18

SAMPLE RESULTS

Lab ID: L1809583-03 Date Collected: 03/20/18 00:00

Client ID: DUP01_032018 Date Received: 03/20/18
Sample Location: BROOKLYN NEW YORK Field Prep: Not Specifie

Sample Location: BROOKLYN, NEW YORK Field Prep: Not Specified

Sample Depth:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | | | |
|--|--------|-----------|-------|-----|------|-----------------|--|--|--|
| Volatile Organics by GC/MS - Westborough Lab | | | | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | | | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | | | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | | | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | | | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | | | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | | | |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 102 | 70-130 | |
| Toluene-d8 | 105 | 70-130 | |
| 4-Bromofluorobenzene | 105 | 70-130 | |
| Dibromofluoromethane | 97 | 70-130 | |

Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/26/18 10:00

| arameter | Result | Qualifier Units | RL | MDL |
|----------------------------|-------------------|-------------------|--------|--------------|
| olatile Organics by GC/MS | - Westborough Lab | for sample(s): 01 | Batch: | WG1100368-12 |
| Methylene chloride | ND | ug/l | 2.5 | 0.70 |
| 1,1-Dichloroethane | ND | ug/l | 2.5 | 0.70 |
| Chloroform | ND | ug/l | 2.5 | 0.70 |
| Carbon tetrachloride | ND | ug/l | 0.50 | 0.13 |
| 1,2-Dichloropropane | ND | ug/l | 1.0 | 0.14 |
| Dibromochloromethane | ND | ug/l | 0.50 | 0.15 |
| 1,1,2-Trichloroethane | ND | ug/l | 1.5 | 0.50 |
| Tetrachloroethene | ND | ug/l | 0.50 | 0.18 |
| Chlorobenzene | ND | ug/l | 2.5 | 0.70 |
| Trichlorofluoromethane | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dichloroethane | ND | ug/l | 0.50 | 0.13 |
| 1,1,1-Trichloroethane | ND | ug/l | 2.5 | 0.70 |
| Bromodichloromethane | ND | ug/l | 0.50 | 0.19 |
| trans-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.16 |
| cis-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.14 |
| 1,3-Dichloropropene, Total | ND | ug/l | 0.50 | 0.14 |
| 1,1-Dichloropropene | ND | ug/l | 2.5 | 0.70 |
| Bromoform | ND | ug/l | 2.0 | 0.65 |
| 1,1,2,2-Tetrachloroethane | ND | ug/l | 0.50 | 0.17 |
| Benzene | ND | ug/l | 0.50 | 0.16 |
| Toluene | ND | ug/l | 2.5 | 0.70 |
| Ethylbenzene | ND | ug/l | 2.5 | 0.70 |
| Chloromethane | ND | ug/l | 2.5 | 0.70 |
| Bromomethane | ND | ug/l | 2.5 | 0.70 |
| Vinyl chloride | ND | ug/l | 1.0 | 0.07 |
| Chloroethane | ND | ug/l | 2.5 | 0.70 |
| 1,1-Dichloroethene | ND | ug/l | 0.50 | 0.17 |
| trans-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 |
| Trichloroethene | ND | ug/l | 0.50 | 0.18 |



Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/26/18 10:00

| Parameter | Result | Qualifier Units | RL | MDL |
|----------------------------|-------------------|-------------------|--------|--------------|
| /olatile Organics by GC/MS | - Westborough Lab | for sample(s): 01 | Batch: | WG1100368-12 |
| 1,2-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,3-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,4-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| Methyl tert butyl ether | ND | ug/l | 2.5 | 0.70 |
| p/m-Xylene | ND | ug/l | 2.5 | 0.70 |
| o-Xylene | ND | ug/l | 2.5 | 0.70 |
| Xylenes, Total | ND | ug/l | 2.5 | 0.70 |
| cis-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dichloroethene, Total | ND | ug/l | 2.5 | 0.70 |
| Dibromomethane | ND | ug/l | 5.0 | 1.0 |
| 1,2,3-Trichloropropane | ND | ug/l | 2.5 | 0.70 |
| Acrylonitrile | ND | ug/l | 5.0 | 1.5 |
| Styrene | ND | ug/l | 2.5 | 0.70 |
| Dichlorodifluoromethane | ND | ug/l | 5.0 | 1.0 |
| Acetone | ND | ug/l | 5.0 | 1.5 |
| Carbon disulfide | ND | ug/l | 5.0 | 1.0 |
| 2-Butanone | ND | ug/l | 5.0 | 1.9 |
| Vinyl acetate | ND | ug/l | 5.0 | 1.0 |
| 4-Methyl-2-pentanone | ND | ug/l | 5.0 | 1.0 |
| 2-Hexanone | ND | ug/l | 5.0 | 1.0 |
| Bromochloromethane | ND | ug/l | 2.5 | 0.70 |
| 2,2-Dichloropropane | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dibromoethane | ND | ug/l | 2.0 | 0.65 |
| 1,3-Dichloropropane | ND | ug/l | 2.5 | 0.70 |
| 1,1,1,2-Tetrachloroethane | ND | ug/l | 2.5 | 0.70 |
| Bromobenzene | ND | ug/l | 2.5 | 0.70 |
| n-Butylbenzene | ND | ug/l | 2.5 | 0.70 |
| sec-Butylbenzene | ND | ug/l | 2.5 | 0.70 |
| tert-Butylbenzene | ND | ug/l | 2.5 | 0.70 |



Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/26/18 10:00

| arameter | Result | Qualifier Units | RL | MDL |
|-------------------------------|---------------|---------------------|--------|--------------|
| olatile Organics by GC/MS - W | estborough La | b for sample(s): 01 | Batch: | WG1100368-12 |
| o-Chlorotoluene | ND | ug/l | 2.5 | 0.70 |
| p-Chlorotoluene | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dibromo-3-chloropropane | ND | ug/l | 2.5 | 0.70 |
| Hexachlorobutadiene | ND | ug/l | 2.5 | 0.70 |
| Isopropylbenzene | ND | ug/l | 2.5 | 0.70 |
| p-Isopropyltoluene | ND | ug/l | 2.5 | 0.70 |
| Naphthalene | ND | ug/l | 2.5 | 0.70 |
| n-Propylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,3-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,3,5-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,4-Dioxane | ND | ug/l | 250 | 61. |
| p-Diethylbenzene | ND | ug/l | 2.0 | 0.70 |
| p-Ethyltoluene | ND | ug/l | 2.0 | 0.70 |
| 1,2,4,5-Tetramethylbenzene | ND | ug/l | 2.0 | 0.54 |
| Ethyl ether | ND | ug/l | 2.5 | 0.70 |
| trans-1,4-Dichloro-2-butene | ND | ug/l | 2.5 | 0.70 |

| | Acceptance | | | | |
|-----------------------|------------|--------------------|--|--|--|
| Surrogate | %Recovery | Qualifier Criteria | | | |
| | | | | | |
| 1,2-Dichloroethane-d4 | 103 | 70-130 | | | |
| Toluene-d8 | 105 | 70-130 | | | |
| 4-Bromofluorobenzene | 105 | 70-130 | | | |
| Dibromofluoromethane | 97 | 70-130 | | | |



Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/25/18 11:00

| Parameter | Result | Qualifier Units | RL | MDL | |
|----------------------------|-------------------|------------------|-------------|-------------|--|
| olatile Organics by GC/MS | - Westborough Lab | for sample(s): 0 | 1-03 Batch: | WG1100368-5 | |
| Methylene chloride | ND | ug/l | 2.5 | 0.70 | |
| 1,1-Dichloroethane | ND | ug/l | 2.5 | 0.70 | |
| Chloroform | ND | ug/l | 2.5 | 0.70 | |
| Carbon tetrachloride | ND | ug/l | 0.50 | 0.13 | |
| 1,2-Dichloropropane | ND | ug/l | 1.0 | 0.14 | |
| Dibromochloromethane | ND | ug/l | 0.50 | 0.15 | |
| 1,1,2-Trichloroethane | ND | ug/l | 1.5 | 0.50 | |
| Tetrachloroethene | ND | ug/l | 0.50 | 0.18 | |
| Chlorobenzene | ND | ug/l | 2.5 | 0.70 | |
| Trichlorofluoromethane | ND | ug/l | 2.5 | 0.70 | |
| 1,2-Dichloroethane | ND | ug/l | 0.50 | 0.13 | |
| 1,1,1-Trichloroethane | ND | ug/l | 2.5 | 0.70 | |
| Bromodichloromethane | ND | ug/l | 0.50 | 0.19 | |
| trans-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.16 | |
| cis-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.14 | |
| 1,3-Dichloropropene, Total | ND | ug/l | 0.50 | 0.14 | |
| 1,1-Dichloropropene | ND | ug/l | 2.5 | 0.70 | |
| Bromoform | ND | ug/l | 2.0 | 0.65 | |
| 1,1,2,2-Tetrachloroethane | ND | ug/l | 0.50 | 0.17 | |
| Benzene | ND | ug/l | 0.50 | 0.16 | |
| Toluene | ND | ug/l | 2.5 | 0.70 | |
| Ethylbenzene | ND | ug/l | 2.5 | 0.70 | |
| Chloromethane | ND | ug/l | 2.5 | 0.70 | |
| Bromomethane | ND | ug/l | 2.5 | 0.70 | |
| Vinyl chloride | ND | ug/l | 1.0 | 0.07 | |
| Chloroethane | ND | ug/l | 2.5 | 0.70 | |
| 1,1-Dichloroethene | ND | ug/l | 0.50 | 0.17 | |
| trans-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 | |
| Trichloroethene | ND | ug/l | 0.50 | 0.18 | |



Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/25/18 11:00

| Parameter | Result | Qualifier Units | RL | MDL |
|---------------------------|-------------------|----------------------|--------|-------------|
| olatile Organics by GC/MS | · Westborough Lab | for sample(s): 01-03 | Batch: | WG1100368-5 |
| 1,2-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,3-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,4-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| Methyl tert butyl ether | ND | ug/l | 2.5 | 0.70 |
| p/m-Xylene | ND | ug/l | 2.5 | 0.70 |
| o-Xylene | ND | ug/l | 2.5 | 0.70 |
| Xylenes, Total | ND | ug/l | 2.5 | 0.70 |
| cis-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dichloroethene, Total | ND | ug/l | 2.5 | 0.70 |
| Dibromomethane | ND | ug/l | 5.0 | 1.0 |
| 1,2,3-Trichloropropane | ND | ug/l | 2.5 | 0.70 |
| Acrylonitrile | ND | ug/l | 5.0 | 1.5 |
| Styrene | ND | ug/l | 2.5 | 0.70 |
| Dichlorodifluoromethane | ND | ug/l | 5.0 | 1.0 |
| Acetone | ND | ug/l | 5.0 | 1.5 |
| Carbon disulfide | ND | ug/l | 5.0 | 1.0 |
| 2-Butanone | ND | ug/l | 5.0 | 1.9 |
| Vinyl acetate | ND | ug/l | 5.0 | 1.0 |
| 4-Methyl-2-pentanone | ND | ug/l | 5.0 | 1.0 |
| 2-Hexanone | ND | ug/l | 5.0 | 1.0 |
| Bromochloromethane | ND | ug/l | 2.5 | 0.70 |
| 2,2-Dichloropropane | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dibromoethane | ND | ug/l | 2.0 | 0.65 |
| 1,3-Dichloropropane | ND | ug/l | 2.5 | 0.70 |
| 1,1,1,2-Tetrachloroethane | ND | ug/l | 2.5 | 0.70 |
| Bromobenzene | ND | ug/l | 2.5 | 0.70 |
| n-Butylbenzene | ND | ug/l | 2.5 | 0.70 |
| sec-Butylbenzene | ND | ug/l | 2.5 | 0.70 |
| tert-Butylbenzene | ND | ug/l | 2.5 | 0.70 |



Project Number: 170357802

Lab Number: L1809583

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/25/18 11:00

| Parameter | Result | Qualifier Units | s RL | MDL |
|--------------------------------|-----------------|-----------------|--------------|-------------|
| Volatile Organics by GC/MS - V | Vestborough Lab | for sample(s): | 01-03 Batch: | WG1100368-5 |
| o-Chlorotoluene | ND | ug/l | 2.5 | 0.70 |
| p-Chlorotoluene | ND | ug/l | | 0.70 |
| 1,2-Dibromo-3-chloropropane | ND | ug/l | 2.5 | 0.70 |
| Hexachlorobutadiene | ND | ug/l | 2.5 | 0.70 |
| Isopropylbenzene | ND | ug/l | 2.5 | 0.70 |
| p-Isopropyltoluene | ND | ug/l | 2.5 | 0.70 |
| Naphthalene | ND | ug/l | 2.5 | 0.70 |
| n-Propylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,3-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,3,5-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,4-Dioxane | ND | ug/l | 250 | 61. |
| p-Diethylbenzene | ND | ug/l | 2.0 | 0.70 |
| p-Ethyltoluene | ND | ug/l | 2.0 | 0.70 |
| 1,2,4,5-Tetramethylbenzene | ND | ug/l | 2.0 | 0.54 |
| Ethyl ether | ND | ug/l | 2.5 | 0.70 |
| trans-1,4-Dichloro-2-butene | ND | ug/l | 2.5 | 0.70 |

| | | Acceptance | | | |
|-----------------------|-----------|--------------------|--|--|--|
| Surrogate | %Recovery | Qualifier Criteria | | | |
| | | | | | |
| 1,2-Dichloroethane-d4 | 101 | 70-130 | | | |
| Toluene-d8 | 104 | 70-130 | | | |
| 4-Bromofluorobenzene | 104 | 70-130 | | | |
| Dibromofluoromethane | 98 | 70-130 | | | |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | %Recovery Qual Limits | RPD | RPD Qual Limits |
|--|------------------|--------------|-------------------|--------------------------|-----|--------------------|
| Volatile Organics by GC/MS - Westborough | Lab Associated | sample(s): 0 | 1 Batch: WG1 | 100368-10 WG1100368-11 | | |
| Methylene chloride | 83 | | 88 | 70-130 | 6 | 20 |
| 1,1-Dichloroethane | 87 | | 93 | 70-130 | 7 | 20 |
| Chloroform | 89 | | 94 | 70-130 | 5 | 20 |
| Carbon tetrachloride | 85 | | 91 | 63-132 | 7 | 20 |
| 1,2-Dichloropropane | 91 | | 97 | 70-130 | 6 | 20 |
| Dibromochloromethane | 96 | | 100 | 63-130 | 4 | 20 |
| 1,1,2-Trichloroethane | 97 | | 100 | 70-130 | 3 | 20 |
| Tetrachloroethene | 95 | | 99 | 70-130 | 4 | 20 |
| Chlorobenzene | 97 | | 100 | 75-130 | 3 | 20 |
| Trichlorofluoromethane | 78 | | 86 | 62-150 | 10 | 20 |
| 1,2-Dichloroethane | 89 | | 96 | 70-130 | 8 | 20 |
| 1,1,1-Trichloroethane | 86 | | 94 | 67-130 | 9 | 20 |
| Bromodichloromethane | 92 | | 99 | 67-130 | 7 | 20 |
| trans-1,3-Dichloropropene | 98 | | 100 | 70-130 | 2 | 20 |
| cis-1,3-Dichloropropene | 92 | | 98 | 70-130 | 6 | 20 |
| 1,1-Dichloropropene | 88 | | 94 | 70-130 | 7 | 20 |
| Bromoform | 88 | | 96 | 54-136 | 9 | 20 |
| 1,1,2,2-Tetrachloroethane | 98 | | 110 | 67-130 | 12 | 20 |
| Benzene | 87 | | 92 | 70-130 | 6 | 20 |
| Toluene | 95 | | 100 | 70-130 | 5 | 20 |
| Ethylbenzene | 97 | | 100 | 70-130 | 3 | 20 |
| Chloromethane | 64 | | 68 | 64-130 | 6 | 20 |
| Bromomethane | 86 | | 88 | 39-139 | 2 | 20 |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|--|------------------|--------------|-------------------|-----------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough L | ab Associated | sample(s): 0 | 1 Batch: WG1 | 100368-10 | WG1100368-11 | | | |
| Vinyl chloride | 76 | | 84 | | 55-140 | 10 | | 20 |
| Chloroethane | 86 | | 92 | | 55-138 | 7 | | 20 |
| 1,1-Dichloroethene | 78 | | 85 | | 61-145 | 9 | | 20 |
| trans-1,2-Dichloroethene | 83 | | 89 | | 70-130 | 7 | | 20 |
| Trichloroethene | 87 | | 92 | | 70-130 | 6 | | 20 |
| 1,2-Dichlorobenzene | 100 | | 100 | | 70-130 | 0 | | 20 |
| 1,3-Dichlorobenzene | 100 | | 100 | | 70-130 | 0 | | 20 |
| 1,4-Dichlorobenzene | 100 | | 100 | | 70-130 | 0 | | 20 |
| Methyl tert butyl ether | 85 | | 94 | | 63-130 | 10 | | 20 |
| p/m-Xylene | 95 | | 100 | | 70-130 | 5 | | 20 |
| o-Xylene | 95 | | 100 | | 70-130 | 5 | | 20 |
| cis-1,2-Dichloroethene | 86 | | 92 | | 70-130 | 7 | | 20 |
| Dibromomethane | 90 | | 99 | | 70-130 | 10 | | 20 |
| 1,2,3-Trichloropropane | 100 | | 110 | | 64-130 | 10 | | 20 |
| Acrylonitrile | 86 | | 95 | | 70-130 | 10 | | 20 |
| Styrene | 95 | | 100 | | 70-130 | 5 | | 20 |
| Dichlorodifluoromethane | 70 | | 79 | | 36-147 | 12 | | 20 |
| Acetone | 90 | | 94 | | 58-148 | 4 | | 20 |
| Carbon disulfide | 77 | | 81 | | 51-130 | 5 | | 20 |
| 2-Butanone | 80 | | 90 | | 63-138 | 12 | | 20 |
| Vinyl acetate | 82 | | 89 | | 70-130 | 8 | | 20 |
| 4-Methyl-2-pentanone | 93 | | 100 | | 59-130 | 7 | | 20 |
| 2-Hexanone | 97 | | 110 | | 57-130 | 13 | | 20 |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | | covery mits RPD | | PD nits |
|---|------------------|--------------|-------------------|----------------|--------------------|-----|------------|
| Volatile Organics by GC/MS - Westboroug | h Lab Associated | sample(s): 0 | 1 Batch: WG1 | 100368-10 WG11 | 00368-11 | | |
| Bromochloromethane | 88 | | 94 | 70 | -130 7 | : | 20 |
| 2,2-Dichloropropane | 89 | | 94 | 63 | -133 5 | | 20 |
| 1,2-Dibromoethane | 97 | | 110 | 70 | -130 13 | | 20 |
| 1,3-Dichloropropane | 97 | | 100 | 70 | -130 3 | | 20 |
| 1,1,1,2-Tetrachloroethane | 96 | | 100 | 64 | -130 4 | | 20 |
| Bromobenzene | 99 | | 100 | 70 | -130 1 | | 20 |
| n-Butylbenzene | 110 | | 110 | 53 | -136 0 | : | 20 |
| sec-Butylbenzene | 110 | | 100 | 70 | -130 10 | - | 20 |
| tert-Butylbenzene | 110 | | 100 | 70 | -130 10 | : | 20 |
| o-Chlorotoluene | 98 | | 99 | 70 | -130 1 | - | 20 |
| p-Chlorotoluene | 100 | | 100 | 70 | -130 0 | - : | 20 |
| 1,2-Dibromo-3-chloropropane | 93 | | 100 | 41 | -144 7 | | 20 |
| Hexachlorobutadiene | 120 | | 120 | 63 | -130 0 | : | 20 |
| Isopropylbenzene | 100 | | 100 | 70 | -130 0 | : | 20 |
| p-lsopropyltoluene | 110 | | 100 | 70 | -130 10 | | 20 |
| Naphthalene | 99 | | 110 | 70 | -130 11 | | 20 |
| n-Propylbenzene | 100 | | 100 | 69 | -130 0 | | 20 |
| 1,2,3-Trichlorobenzene | 100 | | 110 | 70 | -130 10 | : | 20 |
| 1,2,4-Trichlorobenzene | 100 | | 110 | 70 | -130 10 | : | 20 |
| 1,3,5-Trimethylbenzene | 100 | | 100 | 64 | -130 0 | : | 20 |
| 1,2,4-Trimethylbenzene | 100 | | 100 | 70 | -130 0 | : | 20 |
| 1,4-Dioxane | 136 | | 144 | 56 | -162 6 | : | 20 |
| p-Diethylbenzene | 110 | | 110 | 70 | -130 0 | : | 20 |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number:

L1809583

Report Date:

03/28/18

| <u>Parameter</u> | LCS %Recovery | Qual | LCSD %Recovery | ' Qual | %Recovery Limits | RPD | Qual | RPD Limits | |
|---|------------------|--------------|-------------------|-------------|---------------------|-----|------|---------------|--|
| Volatile Organics by GC/MS - Westborough La | ab Associated | sample(s): 0 |)1 Batch: W | G1100368-10 | WG1100368-11 | | | | |
| p-Ethyltoluene | 100 | | 100 | | 70-130 | 0 | | 20 | |
| 1,2,4,5-Tetramethylbenzene | 100 | | 100 | | 70-130 | 0 | | 20 | |
| Ethyl ether | 82 | | 90 | | 59-134 | 9 | | 20 | |
| trans-1,4-Dichloro-2-butene | 94 | | 100 | | 70-130 | 6 | | 20 | |

| | LCS | LCSD | Acceptance |
|-----------------------|----------------|----------------|------------|
| Surrogate | %Recovery Qual | %Recovery Qual | Criteria |
| 1,2-Dichloroethane-d4 | 103 | 104 | 70-130 |
| Toluene-d8 | 104 | 103 | 70-130 |
| 4-Bromofluorobenzene | 102 | 102 | 70-130 |
| Dibromofluoromethane | 99 | 99 | 70-130 |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | %Recovery Qual Limits | RPD | RPD Qual Limits |
|--|------------------|------------|-------------------|--------------------------|-----|--------------------|
| Volatile Organics by GC/MS - Westborough | Lab Associated | sample(s): | 01-03 Batch: W | G1100368-3 WG1100368-4 | | |
| Methylene chloride | 89 | | 86 | 70-130 | 3 | 20 |
| 1,1-Dichloroethane | 93 | | 89 | 70-130 | 4 | 20 |
| Chloroform | 95 | | 91 | 70-130 | 4 | 20 |
| Carbon tetrachloride | 86 | | 84 | 63-132 | 2 | 20 |
| 1,2-Dichloropropane | 97 | | 94 | 70-130 | 3 | 20 |
| Dibromochloromethane | 100 | | 99 | 63-130 | 1 | 20 |
| 1,1,2-Trichloroethane | 100 | | 98 | 70-130 | 2 | 20 |
| Tetrachloroethene | 97 | | 94 | 70-130 | 3 | 20 |
| Chlorobenzene | 100 | | 98 | 75-130 | 2 | 20 |
| Trichlorofluoromethane | 77 | | 76 | 62-150 | 1 | 20 |
| 1,2-Dichloroethane | 93 | | 91 | 70-130 | 2 | 20 |
| 1,1,1-Trichloroethane | 90 | | 87 | 67-130 | 3 | 20 |
| Bromodichloromethane | 96 | | 92 | 67-130 | 4 | 20 |
| trans-1,3-Dichloropropene | 100 | | 100 | 70-130 | 0 | 20 |
| cis-1,3-Dichloropropene | 98 | | 94 | 70-130 | 4 | 20 |
| 1,1-Dichloropropene | 89 | | 87 | 70-130 | 2 | 20 |
| Bromoform | 95 | | 90 | 54-136 | 5 | 20 |
| 1,1,2,2-Tetrachloroethane | 100 | | 100 | 67-130 | 0 | 20 |
| Benzene | 93 | | 89 | 70-130 | 4 | 20 |
| Toluene | 100 | | 97 | 70-130 | 3 | 20 |
| Ethylbenzene | 100 | | 97 | 70-130 | 3 | 20 |
| Chloromethane | 71 | | 69 | 64-130 | 3 | 20 |
| Bromomethane | 88 | | 90 | 39-139 | 2 | 20 |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | RPD Qual Limits | |
|--|------------------|------------|-------------------|-------------|---------------------|-----|--------------------|--|
| Volatile Organics by GC/MS - Westborough I | _ab Associated | sample(s): | 01-03 Batch: W | /G1100368-3 | WG1100368-4 | | | |
| Vinyl chloride | 79 | | 78 | | 55-140 | 1 | 20 | |
| Chloroethane | 90 | | 88 | | 55-138 | 2 | 20 | |
| 1,1-Dichloroethene | 80 | | 78 | | 61-145 | 3 | 20 | |
| trans-1,2-Dichloroethene | 88 | | 84 | | 70-130 | 5 | 20 | |
| Trichloroethene | 92 | | 88 | | 70-130 | 4 | 20 | |
| 1,2-Dichlorobenzene | 100 | | 100 | | 70-130 | 0 | 20 | |
| 1,3-Dichlorobenzene | 110 | | 100 | | 70-130 | 10 | 20 | |
| 1,4-Dichlorobenzene | 100 | | 100 | | 70-130 | 0 | 20 | |
| Methyl tert butyl ether | 90 | | 88 | | 63-130 | 2 | 20 | |
| p/m-Xylene | 100 | | 95 | | 70-130 | 5 | 20 | |
| o-Xylene | 100 | | 95 | | 70-130 | 5 | 20 | |
| cis-1,2-Dichloroethene | 93 | | 89 | | 70-130 | 4 | 20 | |
| Dibromomethane | 94 | | 92 | | 70-130 | 2 | 20 | |
| 1,2,3-Trichloropropane | 100 | | 100 | | 64-130 | 0 | 20 | |
| Acrylonitrile | 91 | | 83 | | 70-130 | 9 | 20 | |
| Styrene | 100 | | 95 | | 70-130 | 5 | 20 | |
| Dichlorodifluoromethane | 69 | | 70 | | 36-147 | 1 | 20 | |
| Acetone | 100 | | 85 | | 58-148 | 16 | 20 | |
| Carbon disulfide | 82 | | 76 | | 51-130 | 8 | 20 | |
| 2-Butanone | 88 | | 76 | | 63-138 | 15 | 20 | |
| Vinyl acetate | 86 | | 83 | | 70-130 | 4 | 20 | |
| 4-Methyl-2-pentanone | 96 | | 91 | | 59-130 | 5 | 20 | |
| 2-Hexanone | 100 | | 97 | | 57-130 | 3 | 20 | |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number: L1809583

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | %Recovery Qual Limits | RPD | RPD Qual Limits |
|--|--------------------|------------|-------------------|--------------------------|-----|--------------------|
| Volatile Organics by GC/MS - Westborou | igh Lab Associated | sample(s): | 01-03 Batch: W | G1100368-3 WG1100368-4 | | |
| Bromochloromethane | 92 | | 89 | 70-130 | 3 | 20 |
| 2,2-Dichloropropane | 94 | | 90 | 63-133 | 4 | 20 |
| 1,2-Dibromoethane | 100 | | 99 | 70-130 | 1 | 20 |
| 1,3-Dichloropropane | 100 | | 99 | 70-130 | 1 | 20 |
| 1,1,1,2-Tetrachloroethane | 100 | | 99 | 64-130 | 1 | 20 |
| Bromobenzene | 100 | | 100 | 70-130 | 0 | 20 |
| n-Butylbenzene | 120 | | 100 | 53-136 | 18 | 20 |
| sec-Butylbenzene | 110 | | 100 | 70-130 | 10 | 20 |
| tert-Butylbenzene | 110 | | 100 | 70-130 | 10 | 20 |
| o-Chlorotoluene | 110 | | 98 | 70-130 | 12 | 20 |
| p-Chlorotoluene | 110 | | 100 | 70-130 | 10 | 20 |
| 1,2-Dibromo-3-chloropropane | 99 | | 95 | 41-144 | 4 | 20 |
| Hexachlorobutadiene | 120 | | 120 | 63-130 | 0 | 20 |
| Isopropylbenzene | 110 | | 100 | 70-130 | 10 | 20 |
| p-Isopropyltoluene | 110 | | 100 | 70-130 | 10 | 20 |
| Naphthalene | 100 | | 99 | 70-130 | 1 | 20 |
| n-Propylbenzene | 110 | | 100 | 69-130 | 10 | 20 |
| 1,2,3-Trichlorobenzene | 100 | | 100 | 70-130 | 0 | 20 |
| 1,2,4-Trichlorobenzene | 110 | | 100 | 70-130 | 10 | 20 |
| 1,3,5-Trimethylbenzene | 110 | | 99 | 64-130 | 11 | 20 |
| 1,2,4-Trimethylbenzene | 110 | | 100 | 70-130 | 10 | 20 |
| 1,4-Dioxane | 138 | | 144 | 56-162 | 4 | 20 |
| p-Diethylbenzene | 120 | | 100 | 70-130 | 18 | 20 |



Project Name: 441 KEAP STREET

Project Number: 170357802

Lab Number:

L1809583

03/28/18

Report Date:

| Parameter | LCS %Recovery | Qual | LCSD %Recover | | %Recovery Limits | RPD | Qual | RPD Limits | |
|--|------------------|------------|------------------|-------------|---------------------|-----|------|---------------|--|
| Volatile Organics by GC/MS - Westborough L | ab Associated | sample(s): | 01-03 Batch: | WG1100368-3 | WG1100368-4 | | | | |
| p-Ethyltoluene | 110 | | 100 | | 70-130 | 10 | | 20 | |
| 1,2,4,5-Tetramethylbenzene | 110 | | 100 | | 70-130 | 10 | | 20 | |
| Ethyl ether | 87 | | 83 | | 59-134 | 5 | | 20 | |
| trans-1,4-Dichloro-2-butene | 100 | | 92 | | 70-130 | 8 | | 20 | |

| | LCS | LCSD | Acceptance | |
|-----------------------|---------------|------------------|------------|---|
| Surrogate | %Recovery Qua | l %Recovery Qual | Criteria | _ |
| 1,2-Dichloroethane-d4 | 102 | 101 | 70-130 | |
| Toluene-d8 | 104 | 105 | 70-130 | |
| 4-Bromofluorobenzene | 101 | 101 | 70-130 | |
| Dibromofluoromethane | 99 | 99 | 70-130 | |

Project Name: 441 KEAP STREET

Lab Number: L1809583

Project Number: 170357802 **Report Date:** 03/28/18

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal

A Absent

| Container Info | ormation | | Initial | Final | Temp | | | Frozen | |
|----------------|--------------------|--------|---------|-------|-------|------|--------|-----------|----------------|
| Container ID | Container Type | Cooler | рН | рН | deg C | Pres | Seal | Date/Time | Analysis(*) |
| L1809583-01A | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-01B | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-01C | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-02A | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-02B | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-02C | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-03A | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-03B | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809583-03C | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |



Project Name: 441 KEAP STREET Lab Number: L1809583

Project Number: 170357802 **Report Date:** 03/28/18

GLOSSARY

Acronyms

EDL - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated

values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis

of PAHs using Solid-Phase Microextraction (SPME).

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LCSD - Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

MDL - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any

adjustments from dilutions, concentrations or moisture content, where applicable.

MS - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for

which an independent estimate of target analyte concentration is available.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less

precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound

list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Footnotes

- The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

Data Qualifiers

A - Spectra identified as "Aldol Condensation Product".

B - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related

Report Format: DU Report with 'J' Qualifiers



Project Name:441 KEAP STREETLab Number:L1809583Project Number:170357802Report Date:03/28/18

Data Qualifiers

projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).

- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations
 of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- M Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.

Report Format: DU Report with 'J' Qualifiers



Project Name:441 KEAP STREETLab Number:L1809583Project Number:170357802Report Date:03/28/18

REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - IV, 2007.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc.
Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:**17873** Revision 11

Published Date: 1/8/2018 4:15:49 PM

Page 1 of 1

Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624: m/p-xylene, o-xylene

EPA 8260C: <u>NPW</u>: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; <u>SCM</u>: lodomethane (methyl iodide), Methyl methacrylate, 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.

EPA 8270D: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

EPA 300: <u>DW:</u> Bromide EPA 6860: <u>SCM:</u> Perchlorate

EPA 9010: NPW and SCM: Amenable Cyanide Distillation

SM4500: NPW: Amenable Cyanide, Dissolved Oxygen; SCM: Total Phosphorus, TKN, NO2, NO3.

Mansfield Facility

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

Drinking Water

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500CN-CE, EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, EPA 351.1, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D.

EPA 624: Volatile Halocarbons & Aromatics,

EPA 608: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, SM9222D.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Be, Cd, Cr, Cu, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn.

EPA 245.1 Hg.

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Document Type: Form Pre-Qualtrax Document ID: 08-113

| ΔLPHA | NEW YORK CHAIN OF CUSTODY | Service Centers Mahwah, NJ 07430; 35 Whitne Albany, NY 12205: 14 Walker V Tonawanda, NY 14150: 275 Co | Vay | 05 | Pag | e i | D | ate Rec'd in Lab | 31 | 2/18 | | ALPHA JOB# L1809583 | |
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| Westbarough, MA 01581 8 Walkup Dr. | Mansfield, MA 02048 320 Forbes Blvd | Project Information | | | *** | | Deliver | | | | | Billing Information | |
| TEL: 508-898-9220 FAX: 508-898-9193 | TEL: 508-822-9300 FAX: 508-822-3288 | | Keup. | | | | - 550 | SP-A | means a | ASP- | | Same as Client Info | |
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| Preservative Code: A = None B = HCI C = HNO ₃ | Container Code P = Plastic A = Amber Glass V = Vial | Westboro: Certification N Mansfield: Certification N | | | | tainer Type | | | | | | Please print clearly, legibly and completely. Samples of not be logged in and | * |
| D = H ₂ SO ₄ E = NaOH | G = Glass B = Bacteria Cup | | 1 | | P | reservative | | | | | | turneround time clock will r | |
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| Westborough, MA 01581 8 Walkup Dr. | 320 Forbes Blvd | Project Information | - | | | | Deliv | erables | | | | Billing Information |
| TEL: 508-898-9220 FAX: 508-898-9193 | TEL: 508-822-9300 FAX: 508-822-3288 | Project Name: 441 | | Street | | | | ASP-A | | X | | Same as Client Info |
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| Preservative Code: | Container Code | Westboro: Certification | L- MAGOS | | | | | | | | | |
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| $F = MeOH$ $G = NaHSO_4$ $H = Na_2S_2O_3$ $K/E = Zn Ac/NaOH$ $O = Other$ Form No: 01-25 HC (rev. 3 | O = Other E = Encore D = BOD Bottle | # Allinguished | 121 | 3/20/18 3/20/18 0/70/2 | | Komeik | hik. | ed By: | 2 | 3/20 | 1707 Serice | resolved. BY EXECUTING THIS COC, THE CLIENT HAS READ AND AGREES TO BE BOUND BY ALPHA'S TERMS & CONDITIONS. (See reverse side.) |



ANALYTICAL REPORT

Lab Number: L1809584

Client: Langan Engineering & Environmental

21 Penn Plaza

360 W. 31st Street, 8th Floor New York, NY 10001-2727

Brian Gochenaur

Phone: (212) 479-5590

Project Name: 123 HOPE STREET

Project Number: 170357801

Report Date: 03/28/18

ATTN:

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), NJ NELAP (MA935), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-14-00197).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number:

L1809584

| Alpha Sample ID | Client ID | Matrix | Sample Location | Collection Date/Time | Receive Date |
|--------------------|--------------|--------|--------------------|----------------------|--------------|
| L1809584-01 | PMW01_032018 | WATER | BROOKLYN, NEW YORK | 03/20/18 10:50 | 03/20/18 |
| L1809584-02 | TB01_032018 | WATER | BROOKLYN, NEW YORK | 03/20/18 11:11 | 03/20/18 |
| L1809584-03 | FB01_032018 | WATER | BROOKLYN, NEW YORK | 03/20/18 11:11 | 03/20/18 |



L1809584

Lab Number:

Project Name: 123 HOPE STREET

Project Number: 170357801 **Report Date:** 03/28/18

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. All specific QC information is also incorporated in the Data Usability format of our Data Merger tool where it can be reviewed along with any associated usability implications. Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances the specific failure is not narrated but noted in the associated QC table. The information is also incorporated in the Data Usability format of our Data Merger tool where it can be reviewed along with any associated usability implications.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

HOLD POLICY

For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Client Service Representative and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

| Please contact (| Client Services at | 800-624-9220 v | with anv c | uestions. |
|------------------|--------------------|----------------|------------|-----------|
| | | | | |



Project Name: 123 HOPE STREET Lab Number: L1809584

Project Number: 170357801 **Report Date:** 03/28/18

Case Narrative (continued)

Report Revision

March 28, 2018: At the client's request, the results of the sample "DUP01_032018" was removed from this report and have been issued under separate cover.

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Amita Naik

Authorized Signature:

Title: Technical Director/Representative Date: 03/28/18

Nails

ORGANICS



VOLATILES



Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1809584

Report Date: 03/28/18

Lab ID: L1809584-01 Date Collected: 03/20/18 10:50

PMW01_032018 Client ID: Date Received: 03/20/18 Sample Location: Field Prep: BROOKLYN, NEW YORK Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/25/18 15:12

Analyst: ΑD

| | | Qualifier | Units | RL | MDL | Dilution Factor |
|---|-------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | h Lab | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | 3.3 | | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |



L1809584

03/28/18

Project Name: 123 HOPE STREET

L1809584-01

Project Number: 170357801

SAMPLE RESULTS

Date Collected: 03/20/18 10:50

Lab Number:

Report Date:

Date Received: 03/20/18

PMW01_032018 Sample Location: Field Prep: Not Specified BROOKLYN, NEW YORK

Sample Depth:

Lab ID:

Client ID:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | gh Lab | | | | | |
| Trichloroethene | 2.0 | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | 23 | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | 23 | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: 123 HOPE STREET **Lab Number:** L1809584

Project Number: 170357801 **Report Date:** 03/28/18

SAMPLE RESULTS

Lab ID: L1809584-01 Date Collected: 03/20/18 10:50

Client ID: PMW01_032018 Date Received: 03/20/18
Sample Location: BROOKLYN, NEW YORK Field Prep: Not Specified

| Volatile Organics by GC/MS - Westborough Lab n-Propylbenzene ND ug/l 2.5 0.70 1 1,2,3-Trichlorobenzene ND ug/l 2.5 0.70 1 1,2,4-Trichlorobenzene ND ug/l 2.5 0.70 1 1,3,5-Trimethylbenzene ND ug/l 2.5 0.70 1 1,2,4-Trimethylbenzene ND ug/l 2.5 0.70 1 1,4-Dioxane ND ug/l 250 61. 1 p-Diethylbenzene ND ug/l 2.0 0.70 1 p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 trans-1,4-Dichloro-2-butene ND ug/l 2.5 0.70 1 | Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|---|----------------------------------|--------------|-----------|-------|-----|------|-----------------|--|
| 1,2,3-Trichlorobenzene ND ug/l 2.5 0.70 1 1,2,4-Trichlorobenzene ND ug/l 2.5 0.70 1 1,3,5-Trimethylbenzene ND ug/l 2.5 0.70 1 1,2,4-Trimethylbenzene ND ug/l 2.5 0.70 1 1,4-Dioxane ND ug/l 250 61. 1 p-Diethylbenzene ND ug/l 2.0 0.70 1 p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | Volatile Organics by GC/MS - Wes | tborough Lab | | | | | | |
| 1,2,4-Trichlorobenzene ND ug/l 2.5 0.70 1 1,3,5-Trimethylbenzene ND ug/l 2.5 0.70 1 1,2,4-Trimethylbenzene ND ug/l 2.5 0.70 1 1,4-Dioxane ND ug/l 250 61. 1 p-Diethylbenzene ND ug/l 2.0 0.70 1 p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene ND ug/l 2.5 0.70 1 1,2,4-Trimethylbenzene ND ug/l 2.5 0.70 1 1,4-Dioxane ND ug/l 250 61. 1 p-Diethylbenzene ND ug/l 2.0 0.70 1 p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene ND ug/l 2.5 0.70 1 1,4-Dioxane ND ug/l 250 61. 1 p-Diethylbenzene ND ug/l 2.0 0.70 1 p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane ND ug/l 250 61. 1 p-Diethylbenzene ND ug/l 2.0 0.70 1 p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| p-Diethylbenzene ND ug/l 2.0 0.70 1 p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| p-Ethyltoluene ND ug/l 2.0 0.70 1 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| 1,2,4,5-Tetramethylbenzene ND ug/l 2.0 0.54 1 Ethyl ether ND ug/l 2.5 0.70 1 | p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| Ethyl ether ND ug/l 2.5 0.70 1 | p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| | 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| trans-1,4-Dichloro-2-butene ND ug/l 2.5 0.70 1 | Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| | trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|-----------------------|------------|-----------|------------------------|--|
| 1,2-Dichloroethane-d4 | 101 | | 70-130 | |
| Toluene-d8 | 105 | | 70-130 | |
| 4-Bromofluorobenzene | 105 | | 70-130 | |
| Dibromofluoromethane | 97 | | 70-130 | |



Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1809584

Report Date: 03/28/18

Result

Lab ID: L1809584-02

Client ID: TB01_032018

Sample Location: BROOKLYN, NEW YORK

Sample Depth:

Parameter

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/25/18 14:17

Analyst: AD

| Date Collected: | 03/20/18 11:11 |
|-----------------|----------------|
| Date Received: | 03/20/18 |
| Field Prep: | Not Specified |

MDL

Dilution Factor

| - uramotor | | | | | |
|----------------------------------|---------------|------|------|------|---|
| Volatile Organics by GC/MS - Wes | stborough Lab | | | | |
| Methylene chloride | ND | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | ND | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | ND | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 | 1 |

Qualifier

Units

RL



L1809584

03/28/18

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number:

Report Date:

Lab ID: L1809584-02 Date Collected: 03/20/18 11:11

Client ID: Date Received: 03/20/18 TB01_032018

Sample Location: Field Prep: Not Specified BROOKLYN, NEW YORK

| Volatile Organics by GC/MS - Westborough Lab Viol ughl 0.50 0.18 1 1.2-Olchiorobexone ND ughl 2.5 0.70 1 1.4-Olchiorobexone ND ughl 2.5 0.70 1 1.4-Olchiorobexone ND ughl 2.5 0.70 1 Methyl feet Luyl ether ND ughl 2.5 0.70 1 PmXylene ND ughl 2.5 0.70 1 Vylene, Total ND ughl 2.5 0.70 1 Vylene, Total ND ughl 2.5 0.70 1 Vylene, Total ND ughl 2.5 0.70 1 Jest-(2-Olchoroethene ND ughl 2.5 0.70 1 Jest-(2-Olchoroethene, Total ND ughl 2.5 0.70 1 Dibroordeflere ND ughl 2.5 0.70 1 Als-(2-Olchoroethene ND ughl 2.5 < | Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------------------------------------|----------|-----------|-------|------|------|-----------------|
| 1.2 Dichlorobenzene | Volatile Organics by GC/MS - Westbor | ough Lab | | | | | |
| 1,2-Dichlorobenzene ND ugil 2,5 0,70 1 1,3-Dichlorobenzene ND ugil 2,5 0,70 1 1,3-Dichlorobenzene ND ugil 2,5 0,70 1 Methyl terb tuyl ether ND ugil 2,5 0,70 1 o-Xylene ND ugil 2,5 0,70 1 o-Xylene ND ugil 2,5 0,70 1 dis-1,2-Dichloroethene ND ugil 2,5 0,70 1 1,2-Dichloroethene, Total ND ugil 2,5 0,70 1 Dibromomethane ND ugil 2,5 0,70 1 Actychirline ND ugil 2,5 0,70 1 Actychirline ND ugil 5,0 1,0 1 Actychirline ND ugil 5,0 1,0 1 Actychirline ND ugil 5,0 1,0 1 | Trichloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| 1,3-Dichlorobenzene ND ugl 2,5 0,70 1 1,4-Dichlorobenzene ND ugl 2,5 0,70 1 Methyl tert buryl ether ND ugl 2,5 0,70 1 o-Sylene ND ugl 2,5 0,70 1 o-Sylene ND ugl 2,5 0,70 1 xylenes, Total ND ugl 2,5 0,70 1 1,2-Dichloroethene ND ugl 2,5 0,70 1 1,2-Dichloroethene, Total ND ugl 2,5 0,70 1 1,2-Dichloroethene, Total ND ugl 2,0 1,0 1 1,2-Dichloroethene, Total ND ugl 2,0 1,0 1 1,2-Dichloroethene, Total ND ugl 2,0 1,0 1 2,2-Dichloroethene, Total ND ugl 2,0 1,0 1 Styrene ND ugl 2,0 1,0 1 <td>1,2-Dichlorobenzene</td> <td>ND</td> <td></td> <td></td> <td>2.5</td> <td>0.70</td> <td>1</td> | 1,2-Dichlorobenzene | ND | | | 2.5 | 0.70 | 1 |
| Methyl tert budyl ether ND ug/l 2.5 0.70 1 p/m-Xylene ND ug/l 2.5 0.70 1 o-Xylene ND ug/l 2.5 0.70 1 o-Xylenes ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 1,2-Dichloroethene, Total ND ug/l 2.5 0.70 1 Dibromemsthane ND ug/l 5.0 1.0 1 Acrylonkride ND ug/l 5.0 1.0 1 Acrylonkride ND ug/l 5.0 1.0 1 Styrene ND ug/l 5.0 1.0 1 Styrene ND ug/l 5.0 1.0 1 Obchtorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.0 1 Vilyi acetate | 1,3-Dichlorobenzene | ND | | | 2.5 | 0.70 | 1 |
| ND | 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene ND ug1 2.5 0.70 1 Xylenes, Total ND ug1 2.5 0.70 1 cis-1,2-Dichloroethene, Total ND ug1 2.5 0.70 1 Dibromomethane ND ug1 2.5 0.70 1 Dibromomethane ND ug1 2.5 0.70 1 Acrylontrile ND ug1 2.5 0.70 1 Acrylontrile ND ug1 2.5 0.70 1 Styrene ND ug1 5.0 1.5 1 Acetone ND ug1 5.0 1.5 1 Acetone ND ug1 5.0 1.0 1 Carbon disulfide ND< | Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total ND ug/l 2.5 0.70 1 cis-1,2-Dichlorcethene ND ug/l 2.5 0.70 1 cis-1,2-Dichlorcethene, Total ND ug/l 2.5 0.70 1 Dichloromethane ND ug/l 2.5 0.70 1 L;2-Dichloroptopane ND ug/l 2.5 0.70 1 Acytonitrile ND ug/l 5.0 1.5 1 Syrene ND ug/l 5.0 1.5 1 Dichlorodfluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Vinyl acetate ND ug/l 5.0 1.0 1 | p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene ND ug/l 2.5 0.70 1 1,2-Dichloroethene, Total ND ug/l 2.5 0.70 1 Dibromomethane ND ug/l 5.0 1.0 1 1,2-Trichloropropane ND ug/l 5.0 0.70 1 Acrylontrile ND ug/l 5.0 0.70 1 Styrene ND ug/l 5.0 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Styria acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 4-Hexthyl-2-pentanone ND ug/l 2.5 0.70 1 | o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total ND ug/l 2,5 0,70 1 | Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane ND ug/l 5.0 1.0 1 1.2.3-Trichloropropane ND ug/l 2.5 0.70 1 Acrylonitrile ND ug/l 5.0 1.5 1 Styrene ND ug/l 5.0 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.5 1 Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.0 1 Viryl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 1,3-Dichropropane ND ug/l 2.5 0.70 1 1,1,1,2-T | cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,3-Trichioropropane ND ug/l 2,5 0,70 1 | 1,2-Dichloroethene, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile ND ug/l 5.0 1.5 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.0 1 1-ynyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 1,2-Dibromochlane ND ug/l 2.5 0.70 1 1,1-1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 1,1 | Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Syrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.0 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 8-Methyl-2-pentanone ND ug/l 5.0 1.0 1 8-Pothachoromethane ND ug/l 2.5 0.70 1 1,2-Distromethane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 | 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodiffluoromethane ND ug/l 5.0 1.0 1 1 1 1 1 1 1 1 1 | Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Acetone ND ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 -Butylbenzene ND ug/l 2.5 0.70 1 < | Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone ND ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 <t< td=""><td>Dichlorodifluoromethane</td><td>ND</td><td></td><td>ug/l</td><td>5.0</td><td>1.0</td><td>1</td></t<> | Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone ND ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 | Acetone | ND | | ug/l | 5.0 | 1.5 | 1 |
| Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 <td>Carbon disulfide</td> <td>ND</td> <td></td> <td>ug/l</td> <td>5.0</td> <td>1.0</td> <td>1</td> | Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 c-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 </td <td>2-Butanone</td> <td>ND</td> <td></td> <td>ug/l</td> <td>5.0</td> <td>1.9</td> <td>1</td> | 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 |
| 2-Hexanone ND ug/l 5.0 1.0 1 | Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane ND | 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.0 0.65 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 | 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2-Dibromoethane ND ug/l 2.0 0.65 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropyltoluene ND ug/l 2.5 0.70 1 | Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 c-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropyltenue ND ug/l 2.5 0.70 1 | 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 sopropylbenzene ND ug/l 2.5 0.70 1 ug/l 2.5 0.70 1 ug/l 2.5 0.70 1 | n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene ND ug/l 2.5 0.70 1 | Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| | Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene ND ug/l 2.5 0.70 1 | p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| | Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: Lab Number: 123 HOPE STREET L1809584

Project Number: Report Date: 170357801 03/28/18

SAMPLE RESULTS

Lab ID: L1809584-02 Date Collected: 03/20/18 11:11

Client ID: Date Received: 03/20/18 TB01_032018 Field Prep: Not Specified

Sample Location: BROOKLYN, NEW YORK

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|-----------------------------------|-------------|-----------|-------|-----|------|-----------------|--|
| Volatile Organics by GC/MS - West | borough Lab | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 102 | 70-130 | |
| Toluene-d8 | 105 | 70-130 | |
| 4-Bromofluorobenzene | 103 | 70-130 | |
| Dibromofluoromethane | 97 | 70-130 | |



03/20/18 11:11

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1809584

Report Date: 03/28/18

Lab ID: L1809584-03 Date Collected:

Client ID: Date Received: 03/20/18 FB01_032018 Sample Location: Field Prep: BROOKLYN, NEW YORK Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/25/18 14:45

Analyst: ΑD

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|----------------------------------|--------------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Wes | tborough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Report Date: 03/28/18

Lab ID: L1809584-03 Client ID: FB01_032018

Sample Location: BROOKLYN, NEW YORK Date Received: 03/20/18

03/20/18 11:11

L1809584

Field Prep:

Lab Number:

Date Collected:

Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|----------------------------------|---------------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Wes | stborough Lab | | | | | |
| Trichloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |
| | | | | | | |



Project Name: 123 HOPE STREET Lab Number: L1809584

Project Number: 170357801 **Report Date:** 03/28/18

SAMPLE RESULTS

Lab ID: L1809584-03 Date Collected: 03/20/18 11:11

Client ID: FB01_032018 Date Received: 03/20/18
Sample Location: BROOKLYN, NEW YORK Field Prep: Not Specified

Ditoonerry, NEW Torric

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|---|--------|-----------|-------|-----|------|-----------------|--|
| Volatile Organics by GC/MS - Westboroug | h Lab | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 103 | 70-130 | |
| Toluene-d8 | 105 | 70-130 | |
| 4-Bromofluorobenzene | 103 | 70-130 | |
| Dibromofluoromethane | 96 | 70-130 | |



Project Number: 170357801

Lab Number: L1809584

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/25/18 11:00

Analyst: PD

| Parameter | Result | Qualifier Units | RL | MDL | |
|----------------------------|-------------------|------------------|-------------|-------------|--|
| olatile Organics by GC/MS | - Westborough Lab | for sample(s): 0 | 1-03 Batch: | WG1100368-5 | |
| Methylene chloride | ND | ug/l | 2.5 | 0.70 | |
| 1,1-Dichloroethane | ND | ug/l | 2.5 | 0.70 | |
| Chloroform | ND | ug/l | 2.5 | 0.70 | |
| Carbon tetrachloride | ND | ug/l | 0.50 | 0.13 | |
| 1,2-Dichloropropane | ND | ug/l | 1.0 | 0.14 | |
| Dibromochloromethane | ND | ug/l | 0.50 | 0.15 | |
| 1,1,2-Trichloroethane | ND | ug/l | 1.5 | 0.50 | |
| Tetrachloroethene | ND | ug/l | 0.50 | 0.18 | |
| Chlorobenzene | ND | ug/l | 2.5 | 0.70 | |
| Trichlorofluoromethane | ND | ug/l | 2.5 | 0.70 | |
| 1,2-Dichloroethane | ND | ug/l | 0.50 | 0.13 | |
| 1,1,1-Trichloroethane | ND | ug/l | 2.5 | 0.70 | |
| Bromodichloromethane | ND | ug/l | 0.50 | 0.19 | |
| trans-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.16 | |
| cis-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.14 | |
| 1,3-Dichloropropene, Total | ND | ug/l | 0.50 | 0.14 | |
| 1,1-Dichloropropene | ND | ug/l | 2.5 | 0.70 | |
| Bromoform | ND | ug/l | 2.0 | 0.65 | |
| 1,1,2,2-Tetrachloroethane | ND | ug/l | 0.50 | 0.17 | |
| Benzene | ND | ug/l | 0.50 | 0.16 | |
| Toluene | ND | ug/l | 2.5 | 0.70 | |
| Ethylbenzene | ND | ug/l | 2.5 | 0.70 | |
| Chloromethane | ND | ug/l | 2.5 | 0.70 | |
| Bromomethane | ND | ug/l | 2.5 | 0.70 | |
| Vinyl chloride | ND | ug/l | 1.0 | 0.07 | |
| Chloroethane | ND | ug/l | 2.5 | 0.70 | |
| 1,1-Dichloroethene | ND | ug/l | 0.50 | 0.17 | |
| trans-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 | |
| Trichloroethene | ND | ug/l | 0.50 | 0.18 | |



Project Number: 170357801

Lab Number: L1809584

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/25/18 11:00

Analyst: PD

| Parameter | Result | Qualifier Units | RL | MDL |
|---------------------------|-------------------|----------------------|--------|-------------|
| olatile Organics by GC/MS | · Westborough Lab | for sample(s): 01-03 | Batch: | WG1100368-5 |
| 1,2-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,3-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,4-Dichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| Methyl tert butyl ether | ND | ug/l | 2.5 | 0.70 |
| p/m-Xylene | ND | ug/l | 2.5 | 0.70 |
| o-Xylene | ND | ug/l | 2.5 | 0.70 |
| Xylenes, Total | ND | ug/l | 2.5 | 0.70 |
| cis-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dichloroethene, Total | ND | ug/l | 2.5 | 0.70 |
| Dibromomethane | ND | ug/l | 5.0 | 1.0 |
| 1,2,3-Trichloropropane | ND | ug/l | 2.5 | 0.70 |
| Acrylonitrile | ND | ug/l | 5.0 | 1.5 |
| Styrene | ND | ug/l | 2.5 | 0.70 |
| Dichlorodifluoromethane | ND | ug/l | 5.0 | 1.0 |
| Acetone | ND | ug/l | 5.0 | 1.5 |
| Carbon disulfide | ND | ug/l | 5.0 | 1.0 |
| 2-Butanone | ND | ug/l | 5.0 | 1.9 |
| Vinyl acetate | ND | ug/l | 5.0 | 1.0 |
| 4-Methyl-2-pentanone | ND | ug/l | 5.0 | 1.0 |
| 2-Hexanone | ND | ug/l | 5.0 | 1.0 |
| Bromochloromethane | ND | ug/l | 2.5 | 0.70 |
| 2,2-Dichloropropane | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dibromoethane | ND | ug/l | 2.0 | 0.65 |
| 1,3-Dichloropropane | ND | ug/l | 2.5 | 0.70 |
| 1,1,1,2-Tetrachloroethane | ND | ug/l | 2.5 | 0.70 |
| Bromobenzene | ND | ug/l | 2.5 | 0.70 |
| n-Butylbenzene | ND | ug/l | 2.5 | 0.70 |
| sec-Butylbenzene | ND | ug/l | 2.5 | 0.70 |
| tert-Butylbenzene | ND | ug/l | 2.5 | 0.70 |



Project Number: 170357801

Lab Number: L1809584

Report Date: 03/28/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/25/18 11:00

Analyst: PD

| arameter | Result | Qualifier Units | s RL | MDL |
|-----------------------------|-----------------|------------------|--------------|-------------|
| olatile Organics by GC/MS - | Westborough Lat | o for sample(s): | 01-03 Batch: | WG1100368-5 |
| o-Chlorotoluene | ND | ug/l | 2.5 | 0.70 |
| p-Chlorotoluene | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dibromo-3-chloropropane | ND | ug/l | 2.5 | 0.70 |
| Hexachlorobutadiene | ND | ug/l | 2.5 | 0.70 |
| Isopropylbenzene | ND | ug/l | 2.5 | 0.70 |
| p-Isopropyltoluene | ND | ug/l | 2.5 | 0.70 |
| Naphthalene | ND | ug/l | 2.5 | 0.70 |
| n-Propylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,3-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,3,5-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,4-Dioxane | ND | ug/l | 250 | 61. |
| p-Diethylbenzene | ND | ug/l | 2.0 | 0.70 |
| p-Ethyltoluene | ND | ug/l | 2.0 | 0.70 |
| 1,2,4,5-Tetramethylbenzene | ND | ug/l | 2.0 | 0.54 |
| Ethyl ether | ND | ug/l | 2.5 | 0.70 |
| trans-1,4-Dichloro-2-butene | ND | ug/l | 2.5 | 0.70 |

| | | | Acceptance | |
|-----------------------|-----------|-----------|------------|--|
| Surrogate | %Recovery | Qualifier | Criteria | |
| | | | | |
| 1,2-Dichloroethane-d4 | 101 | | 70-130 | |
| Toluene-d8 | 104 | | 70-130 | |
| 4-Bromofluorobenzene | 104 | | 70-130 | |
| Dibromofluoromethane | 98 | | 70-130 | |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1809584

Report Date: 03/28/18

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | %Recovery Qual Limits | RPD | RPD Qual Limits |
|--|------------------|------------|-------------------|--------------------------|-----|--------------------|
| Volatile Organics by GC/MS - Westborough | Lab Associated | sample(s): | 01-03 Batch: W | G1100368-3 WG1100368-4 | | |
| Methylene chloride | 89 | | 86 | 70-130 | 3 | 20 |
| 1,1-Dichloroethane | 93 | | 89 | 70-130 | 4 | 20 |
| Chloroform | 95 | | 91 | 70-130 | 4 | 20 |
| Carbon tetrachloride | 86 | | 84 | 63-132 | 2 | 20 |
| 1,2-Dichloropropane | 97 | | 94 | 70-130 | 3 | 20 |
| Dibromochloromethane | 100 | | 99 | 63-130 | 1 | 20 |
| 1,1,2-Trichloroethane | 100 | | 98 | 70-130 | 2 | 20 |
| Tetrachloroethene | 97 | | 94 | 70-130 | 3 | 20 |
| Chlorobenzene | 100 | | 98 | 75-130 | 2 | 20 |
| Trichlorofluoromethane | 77 | | 76 | 62-150 | 1 | 20 |
| 1,2-Dichloroethane | 93 | | 91 | 70-130 | 2 | 20 |
| 1,1,1-Trichloroethane | 90 | | 87 | 67-130 | 3 | 20 |
| Bromodichloromethane | 96 | | 92 | 67-130 | 4 | 20 |
| trans-1,3-Dichloropropene | 100 | | 100 | 70-130 | 0 | 20 |
| cis-1,3-Dichloropropene | 98 | | 94 | 70-130 | 4 | 20 |
| 1,1-Dichloropropene | 89 | | 87 | 70-130 | 2 | 20 |
| Bromoform | 95 | | 90 | 54-136 | 5 | 20 |
| 1,1,2,2-Tetrachloroethane | 100 | | 100 | 67-130 | 0 | 20 |
| Benzene | 93 | | 89 | 70-130 | 4 | 20 |
| Toluene | 100 | | 97 | 70-130 | 3 | 20 |
| Ethylbenzene | 100 | | 97 | 70-130 | 3 | 20 |
| Chloromethane | 71 | | 69 | 64-130 | 3 | 20 |
| Bromomethane | 88 | | 90 | 39-139 | 2 | 20 |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1809584

Report Date: 03/28/18

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | %Recovery Qual Limits | RPD | RPD Qual Limits |
|--|------------------|------------|-------------------|--------------------------|-----|--------------------|
| Volatile Organics by GC/MS - Westborough | n Lab Associated | sample(s): | 01-03 Batch: W | G1100368-3 WG1100368-4 | | |
| Vinyl chloride | 79 | | 78 | 55-140 | 1 | 20 |
| Chloroethane | 90 | | 88 | 55-138 | 2 | 20 |
| 1,1-Dichloroethene | 80 | | 78 | 61-145 | 3 | 20 |
| trans-1,2-Dichloroethene | 88 | | 84 | 70-130 | 5 | 20 |
| Trichloroethene | 92 | | 88 | 70-130 | 4 | 20 |
| 1,2-Dichlorobenzene | 100 | | 100 | 70-130 | 0 | 20 |
| 1,3-Dichlorobenzene | 110 | | 100 | 70-130 | 10 | 20 |
| 1,4-Dichlorobenzene | 100 | | 100 | 70-130 | 0 | 20 |
| Methyl tert butyl ether | 90 | | 88 | 63-130 | 2 | 20 |
| p/m-Xylene | 100 | | 95 | 70-130 | 5 | 20 |
| o-Xylene | 100 | | 95 | 70-130 | 5 | 20 |
| cis-1,2-Dichloroethene | 93 | | 89 | 70-130 | 4 | 20 |
| Dibromomethane | 94 | | 92 | 70-130 | 2 | 20 |
| 1,2,3-Trichloropropane | 100 | | 100 | 64-130 | 0 | 20 |
| Acrylonitrile | 91 | | 83 | 70-130 | 9 | 20 |
| Styrene | 100 | | 95 | 70-130 | 5 | 20 |
| Dichlorodifluoromethane | 69 | | 70 | 36-147 | 1 | 20 |
| Acetone | 100 | | 85 | 58-148 | 16 | 20 |
| Carbon disulfide | 82 | | 76 | 51-130 | 8 | 20 |
| 2-Butanone | 88 | | 76 | 63-138 | 15 | 20 |
| Vinyl acetate | 86 | | 83 | 70-130 | 4 | 20 |
| 4-Methyl-2-pentanone | 96 | | 91 | 59-130 | 5 | 20 |
| 2-Hexanone | 100 | | 97 | 57-130 | 3 | 20 |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1809584

Report Date: 03/28/18

| arameter | LCS %Recovery | Qual | LCSD %Recovery | | %Recovery Limits | RPD | RPD Qual Limits | |
|--|------------------|------------|-------------------|-------------|---------------------|-----|--------------------|--|
| olatile Organics by GC/MS - Westboroug | h Lab Associated | sample(s): | 01-03 Batch: | WG1100368-3 | WG1100368-4 | | | |
| Bromochloromethane | 92 | | 89 | | 70-130 | 3 | 20 | |
| 2,2-Dichloropropane | 94 | | 90 | | 63-133 | 4 | 20 | |
| 1,2-Dibromoethane | 100 | | 99 | | 70-130 | 1 | 20 | |
| 1,3-Dichloropropane | 100 | | 99 | | 70-130 | 1 | 20 | |
| 1,1,1,2-Tetrachloroethane | 100 | | 99 | | 64-130 | 1 | 20 | |
| Bromobenzene | 100 | | 100 | | 70-130 | 0 | 20 | |
| n-Butylbenzene | 120 | | 100 | | 53-136 | 18 | 20 | |
| sec-Butylbenzene | 110 | | 100 | | 70-130 | 10 | 20 | |
| tert-Butylbenzene | 110 | | 100 | | 70-130 | 10 | 20 | |
| o-Chlorotoluene | 110 | | 98 | | 70-130 | 12 | 20 | |
| p-Chlorotoluene | 110 | | 100 | | 70-130 | 10 | 20 | |
| 1,2-Dibromo-3-chloropropane | 99 | | 95 | | 41-144 | 4 | 20 | |
| Hexachlorobutadiene | 120 | | 120 | | 63-130 | 0 | 20 | |
| Isopropylbenzene | 110 | | 100 | | 70-130 | 10 | 20 | |
| p-Isopropyltoluene | 110 | | 100 | | 70-130 | 10 | 20 | |
| Naphthalene | 100 | | 99 | | 70-130 | 1 | 20 | |
| n-Propylbenzene | 110 | | 100 | | 69-130 | 10 | 20 | |
| 1,2,3-Trichlorobenzene | 100 | | 100 | | 70-130 | 0 | 20 | |
| 1,2,4-Trichlorobenzene | 110 | | 100 | | 70-130 | 10 | 20 | |
| 1,3,5-Trimethylbenzene | 110 | | 99 | | 64-130 | 11 | 20 | |
| 1,2,4-Trimethylbenzene | 110 | | 100 | | 70-130 | 10 | 20 | |
| 1,4-Dioxane | 138 | | 144 | | 56-162 | 4 | 20 | |
| p-Diethylbenzene | 120 | | 100 | | 70-130 | 18 | 20 | |



Project Name: 123 HOPE STREET

Project Number: 170357801 Lab Number:

L1809584 Report Date: 03/28/18

| arameter | LCS %Recovery | Qual | LCSD Qual %Recovery | | %Recovery Qual Limits | | Qual | RPD Limits |
|---------------------------------|----------------------------|--------------|------------------------|-------------|--------------------------|----|------|---------------|
| olatile Organics by GC/MS - Wes | stborough Lab Associated s | sample(s): 0 | 01-03 Batch: | WG1100368-3 | WG1100368-4 | | | |
| p-Ethyltoluene | 110 | | 100 | | 70-130 | 10 | | 20 |
| 1,2,4,5-Tetramethylbenzene | 110 | | 100 | | 70-130 | 10 | | 20 |
| Ethyl ether | 87 | | 83 | | 59-134 | 5 | | 20 |
| trans-1,4-Dichloro-2-butene | 100 | | 92 | | 70-130 | 8 | | 20 |

| | LCS | LCSD | Acceptance | |
|-----------------------|----------------|----------------|------------|--|
| Surrogate | %Recovery Qual | %Recovery Qual | Criteria | |
| 1,2-Dichloroethane-d4 | 102 | 101 | 70-130 | |
| Toluene-d8 | 104 | 105 | 70-130 | |
| 4-Bromofluorobenzene | 101 | 101 | 70-130 | |
| Dibromofluoromethane | 99 | 99 | 70-130 | |



Serial_No:03281815:44 *Lab Number:* L1809584

Project Name: 123 HOPE STREET

Project Number: 170357801 **Report Date:** 03/28/18

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Cooler Custody Seal

A Absent

| Container Info | rmation | | Initial | Final | Temp | | | Frozen | |
|----------------|--------------------|--------|---------|-------|-------|------|--------|-----------|----------------|
| Container ID | Container Type | Cooler | рН | рН | deg C | Pres | Seal | Date/Time | Analysis(*) |
| L1809584-01A | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-01B | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-01C | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-02A | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-02B | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-02C | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-02D | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-03A | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-03B | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-03C | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | NYTCL-8260(14) |
| L1809584-04A | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | - |
| L1809584-04B | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | - |
| L1809584-04C | Vial HCl preserved | Α | NA | | 5.8 | Υ | Absent | | - |
| | | | | | | | | | |



Project Name:123 HOPE STREETLab Number:L1809584Project Number:170357801Report Date:03/28/18

GLOSSARY

Acronyms

EDL - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated

values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis

of PAHs using Solid-Phase Microextraction (SPME).

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LCSD - Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

MDL - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any

adjustments from dilutions, concentrations or moisture content, where applicable.

MS - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for

which an independent estimate of target analyte concentration is available.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less

than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound

list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Footnotes

- The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

Data Qualifiers

A - Spectra identified as "Aldol Condensation Product".

- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related

Report Format: DU Report with 'J' Qualifiers



В

Project Name:123 HOPE STREETLab Number:L1809584Project Number:170357801Report Date:03/28/18

Data Qualifiers

projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).

- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations
 of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- M Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.

Report Format: DU Report with 'J' Qualifiers



 Project Name:
 123 HOPE STREET
 Lab Number:
 L1809584

 Project Number:
 170357801
 Report Date:
 03/28/18

REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - IV, 2007.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc.
Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:**17873** Revision 11

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Published Date: 1/8/2018 4:15:49 PM

Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624: m/p-xylene, o-xylene

EPA 8260C: <u>NPW</u>: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; <u>SCM</u>: lodomethane (methyl iodide), Methyl methacrylate, 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.

EPA 8270D: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

EPA 300: <u>DW:</u> Bromide EPA 6860: <u>SCM:</u> Perchlorate

EPA 9010: NPW and SCM: Amenable Cyanide Distillation

SM4500: NPW: Amenable Cyanide, Dissolved Oxygen; SCM: Total Phosphorus, TKN, NO2, NO3.

Mansfield Facility

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

Drinking Water

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE, EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, EPA 351.1, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D.

EPA 624: Volatile Halocarbons & Aromatics,

EPA 608: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, SM9222D.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Be, Cd, Cr, Cu, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn.

EPA 245.1 Hg.

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Document Type: Form Pre-Qualtrax Document ID: 08-113

| 100 | NEW YORK | Service Centers | | | _ | | _ | | | | | |
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| Westborough, MA 01581 8 Walkup Dr. | Mansfield, MA 02048 320 Forbes Blvd | Project Information | | THE | | | Deliver | ables | 31 | 7.410 | | Billing Information |
| TEL: 508-898-9220 | TEL: 508-822-9300 | Project Name: 123 | Hope S | tient | | | | SP-A | | X ASF | P-B | Same as Client Info |
| FAX: 508-898-9193 | FAX: 508-822-3288 | Project Location: Big | cklya N | OLD YOU | k | | | QuIS (1 | File) | Annual Property of the Parket | ulS (4 File) | PO# |
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| H ₂ SO ₄ G = Glass | | | | | Р | reservative | | 1 1 | | | | not be logged in and turnaround time clock will no |
| | B = Bacteria Cup C = Cube | 77 | | | | | | | | | | start until any ambiguities are |
| = NaHSO ₄ 0 = Other Bellinguished By Date/Time | | | | | | | Rebeived | | | , Date/ | Time | resolved, BY EXECUTING |
| 11020203 | E = Encore D = 80D Bottle | 190 | - | | 17:10 | Kimek | Arcks. | n Aft | 3 | 20 1 | 707 | THIS COC, THE CLIENT |
| /E = Zn Ac/NaOH = Other | - SOD Dollie | 1 | | | 1260 | / | 53 | 120 | | 772 | 0 | HAS READ AND AGREES TO BE BOUND BY ALPHA'S |
| | | -7 | 3/21 | ou | | Organi | F.0 | wa (A) | 3 | | 02:00 | TERMS & CONDITIONS. |
| orm No: 01-25 HC (rev. 30 | 0-Sept-2013) | | | | | 9 9 | U | | 10 | | | (See reverse side.) |



ANALYTICAL REPORT

Lab Number: L1810069

Client: Langan Engineering & Environmental

21 Penn Plaza

360 W. 31st Street, 8th Floor New York, NY 10001-2727

ATTN: Brian Gochenaur Phone: (212) 479-5590

Project Name: 123 HOPE STREET

Project Number: 170357801 Report Date: 03/29/18

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), NJ NELAP (MA935), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-14-00197).

Eight Walkup Drive, Westborough, MA 01581-1019 508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Number: 170357801

Lab Number: L1810069 **Report Date:** 03/29/18

| Alpha Sample ID | Client ID | Matrix | Sample Location | Collection Date/Time | Receive Date |
|--------------------|--------------|--------|--------------------|----------------------|--------------|
| L1810069-01 | PMW02_032318 | WATER | BROOKLYN, NY | 03/23/18 14:50 | 03/23/18 |
| L1810069-02 | PMW03_032318 | WATER | BROOKLYN, NY | 03/23/18 13:38 | 03/23/18 |
| L1810069-03 | PMW04_032318 | WATER | BROOKLYN, NY | 03/23/18 12:20 | 03/23/18 |
| L1810069-04 | PMW05_032318 | WATER | BROOKLYN, NY | 03/23/18 15:40 | 03/23/18 |
| L1810069-05 | TB02_032318 | WATER | BROOKLYN, NY | 03/23/18 00:00 | 03/23/18 |



L1810069

Lab Number:

Project Name: 123 HOPE STREET

Project Number: 170357801 **Report Date:** 03/29/18

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively. When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. All specific QC information is also incorporated in the Data Usability format of our Data Merger tool where it can be reviewed along with any associated usability implications. Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances the specific failure is not narrated but noted in the associated QC table. The information is also incorporated in the Data Usability format of our Data Merger tool where it can be reviewed along with any associated usability implications.

Please see the associated ADEx data file for a comparison of laboratory reporting limits that were achieved with the regulatory Numerical Standards requested on the Chain of Custody.

HOLD POLICY

For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Client Service Representative and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

| Please | contact | Client | Services | at 800- | -624-9220 | with a | nv c | nuestions |
|--------|---------|--------|-----------|---------|-----------|--------|------|-------------------------|
| loase | Contact | Olicit | OCI VICCO | at ooo | 02- 0220 | with a | ıy c | _f ucsiloris. |



L1810069

Project Name: 123 HOPE STREET

Project Number: 170357801 **Report Date:** 03/29/18

Lab Number:

Case Narrative (continued)

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Michelle M. Morris

Authorized Signature:

Title: Technical Director/Representative

Date: 03/29/18

ORGANICS



VOLATILES



03/23/18 14:50

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1810069

Report Date: 03/29/18

Date Collected:

Lab ID: L1810069-01

Client ID: PMW02_032318 Sample Location: BROOKLYN, NY Date Received: 03/23/18 Field Prep: Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/28/18 12:21

Analyst: NLK

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | ıh Lab | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | 15 | | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | 0.13 | J | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |



03/23/18 14:50

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1810069

Report Date: 03/29/18

Lab ID: L1810069-01 Date Collected:

PMW02_032318

BROOKLYN, NY

Date Received: 03/23/18

Field Prep: Not Specified

Sample Depth:

Sample Location:

Client ID:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|-------------------------------------|-----------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westbo | rough Lab | | | | | |
| Trichloroethene | 3.4 | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | 23 | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | 23 | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | ND | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: 123 HOPE STREET Lab Number: L1810069

Project Number: 170357801 **Report Date:** 03/29/18

SAMPLE RESULTS

Lab ID: L1810069-01 Date Collected: 03/23/18 14:50

Client ID: PMW02_032318 Date Received: 03/23/18 Sample Location: BROOKLYN, NY Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|-----------------------------------|-------------|-----------|-------|-----|------|-----------------|--|
| Volatile Organics by GC/MS - West | borough Lab | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 96 | 70-130 | |
| Toluene-d8 | 106 | 70-130 | |
| 4-Bromofluorobenzene | 113 | 70-130 | |
| Dibromofluoromethane | 93 | 70-130 | |

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1810069

Report Date: 03/29/18

Date Collected:

Lab ID: L1810069-02

Client ID: PMW03_032318 Sample Location: BROOKLYN, NY Date Received: 03/23/18 Field Prep: Not Specified

03/23/18 13:38

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/28/18 12:46

Analyst: NLK

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | | |
|--|--------|-----------|-------|------|------|-----------------|--|--|
| Volatile Organics by GC/MS - Westborough Lab | | | | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Chloroform | 1.3 | J | ug/l | 2.5 | 0.70 | 1 | | |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 | | |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 | | |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 | | |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 | | |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 | 1 | | |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 | | |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 | | |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 | | |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 | | |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 | | |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 | | |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 | | |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 | | |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Bromomethane | 1.7 | J | ug/l | 2.5 | 0.70 | 1 | | |
| Vinyl chloride | 0.77 | J | ug/l | 1.0 | 0.07 | 1 | | |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 | | |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 | | |



L1810069

Project Name: 123 HOPE STREET Lab Number:

Project Number: 170357801 **Report Date:** 03/29/18

SAMPLE RESULTS

Lab ID: L1810069-02 Date Collected: 03/23/18 13:38

Client ID: PMW03_032318 Date Received: 03/23/18 Sample Location: BROOKLYN, NY Field Prep: Not Specified

| 1,2,3-Trichloropropane ND ug/l 2.5 0.70 1 | Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--|--------------------------------|----------------|-----------|-------|------|------|-----------------|
| 1,2-Dichlorobenzene ND ugil 2,5 0,70 1 1,3-Dichlorobenzene ND ugil 2,5 0,70 1 1,4-Dichlorobenzene R-Total ND ugil 2,5 0,70 1 1,4-Dichlorobenzene R-Total R-Tot | Volatile Organics by GC/MS - W | estborough Lab | | | | | |
| 1,2-Dichlorobenzene ND ugil 2,5 0,70 1 1,3-Dichlorobenzene ND ugil 2,5 0,70 1 1,4-Dichlorobenzene R-Total ND ugil 2,5 0,70 1 1,4-Dichlorobenzene R-Total R-Tot | Trichloroothono | ND | | ua/l | 0.50 | 0.18 | 1 |
| 1,3-Dichlorobenzone ND ug/l 2,5 0,70 1 1,4-Bichlorobenzene ND ug/l 2,5 0,70 1 1,4-Bichlorobenzen | | | | | | | |
| 1,4-Dichlorobenzene ND Ug/l 2,5 0,70 1 | | | | | | | |
| Methyl tert butyl other ND ug/l 2.5 0.70 1 prim-Xylene ND ug/l 2.5 0.70 1 o-Xylene ND ug/l 2.5 0.70 1 o-Xylenes ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene 6.7 ug/l 2.5 0.70 1 1,2-Dichloroethene, Total 6.7 ug/l 2.5 0.70 1 Dibromomethane ND ug/l 5.0 1.0 1 Dibromomethane ND ug/l 5.0 1.0 1 1,2,3-Trichloropropane ND ug/l 5.0 1.0 1 Styrene ND ug/l 5.0 1.0 1 Styrene ND ug/l 5.0 1.0 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone 18 ug/l 5.0 1.0 1 Vir | | | | | | | |
| p/m-Xylene ND ug/l 2.5 0.70 1 o-Xylene ND ug/l 2.5 0.70 1 xylenes, Total ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene 6.7 ug/l 2.5 0.70 1 1,2-Dichloroethene, Total 6.7 ug/l 2.5 0.70 1 1,2-Dichloroethene, Total ND ug/l 2.5 0.70 1 1,2-Dichloroethene ND ug/l 2.5 0.70 1 Styrone ND ug/l 2.5 0.70 1 Acetone 18 ug/l 5.0 1.0 1 Styrone ND ug/l 5.0 1.0 1 | | | | | | | |
| o-Xylene ND ug/l 2.5 0.70 1 Xylenes, Total ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene 6.7 ug/l 2.5 0.70 1 1,2-Dichloroethene, Total 6.7 ug/l 2.5 0.70 1 Dibromomethane ND ug/l 5.0 1.0 1 1,2-Dichloroethene, Total ND ug/l 5.0 1.0 1 1,2-S-Trichloropropane ND ug/l 5.0 1.5 1 Actyclontifile ND ug/l 5.0 1.5 1 Styrene ND ug/l 5.0 1.0 1 Actione 18 ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 Carbon disulfide ND ug/l 5.0 1.0 1 | | | | | | | |
| Xylenes, Total ND ug/l 2.5 0.70 1 cis-1,2-Dichloroethene 6.7 ug/l 2.5 0.70 1 1,2-Dichloroethene, Total 6.7 ug/l 2.5 0.70 1 Dibromomethane ND ug/l 2.5 0.70 1 Acrylonitrile ND ug/l 2.5 0.70 1 Styrene ND ug/l 5.0 1.5 1 Styrene ND ug/l 5.0 1.0 1 Dichloroffluoromethane ND ug/l 5.0 1.0 1 Acetone 18 ug/l 5.0 1.0 1 Acetone 18 ug/l 5.0 1.0 1 Viryl acetata ND ug/l 5.0 1.0 1 Viryl acetata ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Pethylacet | <u> </u> | | | | | | |
| Cis-1,2-Dichloroethene 6.7 ug/l 2.5 0.70 1 | | | | | | | |
| 1,2-Dichloroethene, Total 6.7 ug/l 2.5 0.70 1 | · | | | | | | |
| Dibromomethane ND ug/l 5.0 1.0 1 1 1 1 1 1 1 1 1 | | | | | | | |
| 1,2,3-Trichloropropane ND ug/l 2,5 0,70 1 | Dibromomethane | | | | | | 1 |
| Actylonitrile ND ug/l 5.0 1.5 1 Styrene ND ug/l 2.5 0.70 1 Dichlorodiflurormethane ND ug/l 5.0 1.0 1 Acetone 18 ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone 2.3 J ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 2.5 0.70 1 2-Hexanone ND ug/l 2.5 0.70 1 Bromochloromethane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 <td>1,2,3-Trichloropropane</td> <td>ND</td> <td></td> <td></td> <td>2.5</td> <td>0.70</td> <td>1</td> | 1,2,3-Trichloropropane | ND | | | 2.5 | 0.70 | 1 |
| Styrene ND ug/l 2.5 0.70 1 Dichlorodifluoromethane ND ug/l 5.0 1.0 1 Acetone 18 ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone 2.3 J ug/l 5.0 1.0 1 Viryl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2-Polichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,1,1-E-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 | Acrylonitrile | ND | | | 5.0 | 1.5 | 1 |
| ND | Styrene | ND | | | 2.5 | 0.70 | 1 |
| Actione 18 ug/l 5.0 1.5 1 Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone 2.3 J ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 2.5 0.70 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 | Dichlorodifluoromethane | ND | | | 5.0 | 1.0 | 1 |
| Carbon disulfide ND ug/l 5.0 1.0 1 2-Butanone 2.3 J ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 </td <td>Acetone</td> <td>18</td> <td></td> <td></td> <td>5.0</td> <td>1.5</td> <td>1</td> | Acetone | 18 | | | 5.0 | 1.5 | 1 |
| 2-Butanone 2.3 J ug/l 5.0 1.9 1 Vinyl acetate ND ug/l 5.0 1.0 1 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1< | Carbon disulfide | ND | | | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone ND ug/l 5.0 1.0 1 2-Hexanone ND ug/l 5.0 1.0 1 Bromochloromethane ND ug/l 2.5 0.70 1 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.5 0.70 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 | 2-Butanone | 2.3 | J | | 5.0 | 1.9 | 1 |
| ND | Vinyl acetate | ND | | | 5.0 | 1.0 | 1 |
| Bromochloromethane ND | 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2,2-Dichloropropane ND ug/l 2.5 0.70 1 1,2-Dibromoethane ND ug/l 2.0 0.65 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 l-2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 tetr-Butylbenzene ND ug/l 2.5 0.70 1 | 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2-Dibromoethane ND ug/l 2.0 0.65 1 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichloropropane ND ug/l 2.5 0.70 1 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| Bromobenzene ND ug/l 2.5 0.70 1 n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene ND ug/l 2.5 0.70 1 sec-Butylbenzene ND ug/l 2.5 0.70 1 tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| ND | Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene ND ug/l 2.5 0.70 1 o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene ND ug/l 2.5 0.70 1 p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene ND ug/l 2.5 0.70 1 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane ND ug/l 2.5 0.70 1 Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene ND ug/l 2.5 0.70 1 Isopropylbenzene ND ug/l 2.5 0.70 1 p-Isopropyltoluene ND ug/l 2.5 0.70 1 | p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene ND ug/l 2.5 0.70 1 | Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| · | Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene ND ug/l 2.5 0.70 1 | p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| | Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: 123 HOPE STREET Lab Number: L1810069

Project Number: 170357801 **Report Date:** 03/29/18

SAMPLE RESULTS

Lab ID: L1810069-02 Date Collected: 03/23/18 13:38

Client ID: PMW03_032318 Date Received: 03/23/18 Sample Location: BROOKLYN, NY Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|---|--------|-----------|-------|-----|------|-----------------|--|
| Volatile Organics by GC/MS - Westboroug | h Lab | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria | |
|-----------------------|------------|-----------|------------------------|--|
| 1,2-Dichloroethane-d4 | 95 | | 70-130 | |
| Toluene-d8 | 106 | | 70-130 | |
| 4-Bromofluorobenzene | 114 | | 70-130 | |
| Dibromofluoromethane | 91 | | 70-130 | |



Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1810069

Report Date: 03/29/18

Lab ID: L1810069-03 Date Collected: 03/23/18 12:20

Client ID: Date Received: 03/23/18 PMW04_032318 Sample Location: Field Prep: BROOKLYN, NY Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/28/18 13:12

Analyst: NLK

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | | |
|--|--------|-----------|-------|------|------|-----------------|--|--|
| Volatile Organics by GC/MS - Westborough Lab | | | | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 | | |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 | | |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 | | |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 | | |
| Tetrachloroethene | 2.5 | | ug/l | 0.50 | 0.18 | 1 | | |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 | | |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 | | |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 | | |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 | | |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 | | |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 | | |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 | | |
| Benzene | 0.37 | J | ug/l | 0.50 | 0.16 | 1 | | |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| Vinyl chloride | 0.89 | J | ug/l | 1.0 | 0.07 | 1 | | |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 | | |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 | | |
| trans-1,2-Dichloroethene | 0.77 | J | ug/l | 2.5 | 0.70 | 1 | | |



L1810069

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Report Date: 03/29/18

Lab Number:

Lab ID: L1810069-03

PMW04_032318 Client ID: Sample Location: BROOKLYN, NY Date Collected: 03/23/18 12:20

Date Received: 03/23/18 Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--------------------------------------|----------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westbor | ough Lab | | | | | |
| Trichloroethene | 1.6 | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | 39 | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | 40 | J | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | 34 | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | 31 | | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: 123 HOPE STREET Lab Number: L1810069

Project Number: 170357801 **Report Date:** 03/29/18

SAMPLE RESULTS

Lab ID: L1810069-03 Date Collected: 03/23/18 12:20

Client ID: PMW04_032318 Date Received: 03/23/18 Sample Location: BROOKLYN, NY Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|-----------------------------------|-------------|-----------|-------|-----|------|-----------------|--|
| Volatile Organics by GC/MS - West | borough Lab | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 99 | 70-130 | |
| Toluene-d8 | 106 | 70-130 | |
| 4-Bromofluorobenzene | 113 | 70-130 | |
| Dibromofluoromethane | 96 | 70-130 | |



03/23/18 15:40

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1810069

Report Date: 03/29/18

Lab ID: L1810069-04 Date Collected:

PMW05_032318 Client ID: Date Received: 03/23/18 Sample Location: Field Prep: BROOKLYN, NY Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/28/18 13:37

Analyst: NLK

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|----------------------------------|--------------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Wes | tborough Lab | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | 0.48 | J | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| | | | | | | |



Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

L1810069

Lab Number:

Report Date: 03/29/18

Lab ID: L1810069-04

PMW05_032318 Client ID: Sample Location: BROOKLYN, NY Date Collected: 03/23/18 15:40

Date Received: 03/23/18 Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|-----------------------------------|--------------|-----------|--------------|------|------|-----------------|
| Volatile Organics by GC/MS - West | tborough Lab | | | | | |
| Trichloroethene | ND | | /l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | | 5.0 | 1.0 | 1 |
| Acetone | 22 | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | 8.7 | | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l ug/l | 2.5 | 0.03 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | | 2.5 | 0.70 | |
| sec-Butylbenzene | ND | | ug/l ug/l | 2.5 | 0.70 | 1 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | | 2.5 | 0.70 | 1 |
| Naphulalelle | ואט | | ug/l | 2.0 | 0.70 | I |



Project Name: 123 HOPE STREET Lab Number: L1810069

Project Number: 170357801 **Report Date:** 03/29/18

SAMPLE RESULTS

Lab ID: L1810069-04 Date Collected: 03/23/18 15:40

Client ID: PMW05_032318 Date Received: 03/23/18 Sample Location: BROOKLYN, NY Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor | |
|---|--------|-----------|-------|-----|------|-----------------|--|
| Volatile Organics by GC/MS - Westboroug | h Lab | | | | | | |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 | |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 | |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 | |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 | |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 | |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 | |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 96 | 70-130 | |
| Toluene-d8 | 107 | 70-130 | |
| 4-Bromofluorobenzene | 113 | 70-130 | |
| Dibromofluoromethane | 92 | 70-130 | |



Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number: L1810069

Report Date: 03/29/18

Lab ID: L1810069-05 Date Collected: 03/23/18 00:00

Client ID: Date Received: 03/23/18 TB02_032318 Sample Location: Field Prep: BROOKLYN, NY Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 03/28/18 14:02

Analyst: NLK

| | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westboroug | h Lab | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,3-Dichloropropene, Total | ND | | ug/l | 0.50 | 0.14 | 1 |
| 1,1-Dichloropropene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |



L1810069

03/23/18 00:00

Project Name: 123 HOPE STREET

Project Number: 170357801

SAMPLE RESULTS

Lab Number:

Report Date: 03/29/18

Lab ID: L1810069-05 Date Collected:

Client ID: Date Received: 03/23/18 TB02_032318 Sample Location: Field Prep: Not Specified BROOKLYN, NY

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough | Lab | | | | | |
| Trichloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Xylenes, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethene, Total | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dibromomethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| 1,2,3-Trichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Acrylonitrile | ND | | ug/l | 5.0 | 1.5 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | 2.6 | J | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 |
| Vinyl acetate | ND | | ug/l | 5.0 | 1.0 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 2,2-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,3-Dichloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1,1,2-Tetrachloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| n-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| sec-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| tert-Butylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Chlorotoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| p-Isopropyltoluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Naphthalene | ND | | ug/l | 2.5 | 0.70 | 1 |



Project Name: 123 HOPE STREET Lab Number: L1810069

Project Number: 170357801 **Report Date:** 03/29/18

SAMPLE RESULTS

Lab ID: L1810069-05 Date Collected: 03/23/18 00:00

Client ID: TB02_032318 Date Received: 03/23/18 Sample Location: BROOKLYN, NY Field Prep: Not Specified

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--|--------|-----------|-------|-----|------|-----------------|
| Volatile Organics by GC/MS - Westborough | Lab | | | | | |
| | | | | | | , |
| n-Propylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,3,5-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,4-Trimethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 |
| p-Diethylbenzene | ND | | ug/l | 2.0 | 0.70 | 1 |
| p-Ethyltoluene | ND | | ug/l | 2.0 | 0.70 | 1 |
| 1,2,4,5-Tetramethylbenzene | ND | | ug/l | 2.0 | 0.54 | 1 |
| Ethyl ether | ND | | ug/l | 2.5 | 0.70 | 1 |
| trans-1,4-Dichloro-2-butene | ND | | ug/l | 2.5 | 0.70 | 1 |

| Surrogate | % Recovery | Acceptance Qualifier Criteria | |
|-----------------------|------------|----------------------------------|--|
| 1,2-Dichloroethane-d4 | 98 | 70-130 | |
| Toluene-d8 | 107 | 70-130 | |
| 4-Bromofluorobenzene | 114 | 70-130 | |
| Dibromofluoromethane | 92 | 70-130 | |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1810069

03/29/18

Report Date:

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/28/18 09:49

Analyst: PD

| Parameter | Result | Qualifier Units | RL | MDL | |
|----------------------------|-------------------|------------------|-------------|-------------|--|
| olatile Organics by GC/MS | - Westborough Lab | for sample(s): 0 | 1-05 Batch: | WG1101340-5 | |
| Methylene chloride | ND | ug/l | 2.5 | 0.70 | |
| 1,1-Dichloroethane | ND | ug/l | 2.5 | 0.70 | |
| Chloroform | ND | ug/l | 2.5 | 0.70 | |
| Carbon tetrachloride | ND | ug/l | 0.50 | 0.13 | |
| 1,2-Dichloropropane | ND | ug/l | 1.0 | 0.14 | |
| Dibromochloromethane | ND | ug/l | 0.50 | 0.15 | |
| 1,1,2-Trichloroethane | ND | ug/l | 1.5 | 0.50 | |
| Tetrachloroethene | ND | ug/l | 0.50 | 0.18 | |
| Chlorobenzene | ND | ug/l | 2.5 | 0.70 | |
| Trichlorofluoromethane | ND | ug/l | 2.5 | 0.70 | |
| 1,2-Dichloroethane | ND | ug/l | 0.50 | 0.13 | |
| 1,1,1-Trichloroethane | ND | ug/l | 2.5 | 0.70 | |
| Bromodichloromethane | ND | ug/l | 0.50 | 0.19 | |
| trans-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.16 | |
| cis-1,3-Dichloropropene | ND | ug/l | 0.50 | 0.14 | |
| 1,3-Dichloropropene, Total | ND | ug/l | 0.50 | 0.14 | |
| 1,1-Dichloropropene | ND | ug/l | 2.5 | 0.70 | |
| Bromoform | ND | ug/l | 2.0 | 0.65 | |
| 1,1,2,2-Tetrachloroethane | ND | ug/l | 0.50 | 0.17 | |
| Benzene | ND | ug/l | 0.50 | 0.16 | |
| Toluene | ND | ug/l | 2.5 | 0.70 | |
| Ethylbenzene | ND | ug/l | 2.5 | 0.70 | |
| Chloromethane | ND | ug/l | 2.5 | 0.70 | |
| Bromomethane | ND | ug/l | 2.5 | 0.70 | |
| Vinyl chloride | ND | ug/l | 1.0 | 0.07 | |
| Chloroethane | ND | ug/l | 2.5 | 0.70 | |
| 1,1-Dichloroethene | ND | ug/l | 0.50 | 0.17 | |
| trans-1,2-Dichloroethene | ND | ug/l | 2.5 | 0.70 | |
| Trichloroethene | ND | ug/l | 0.50 | 0.18 | |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1810069

Report Date: 03/29/18

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/28/18 09:49

Analyst: PD

| arameter | Result | Qualifier Units | s RL | MDL |
|-----------------------------|----------------|------------------|--------------|-------------|
| olatile Organics by GC/MS - | Westborough La | b for sample(s): | 01-05 Batch: | WG1101340-5 |
| 1,2-Dichlorobenzene | ND | ug/ | l 2.5 | 0.70 |
| 1,3-Dichlorobenzene | ND | ug/ | l 2.5 | 0.70 |
| 1,4-Dichlorobenzene | ND | ug/ | 1 2.5 | 0.70 |
| Methyl tert butyl ether | ND | ug/ | 1 2.5 | 0.70 |
| p/m-Xylene | ND | ug/ | 1 2.5 | 0.70 |
| o-Xylene | ND | ug/ | l 2.5 | 0.70 |
| Xylenes, Total | ND | ug/ | 1 2.5 | 0.70 |
| cis-1,2-Dichloroethene | ND | ug/ | 1 2.5 | 0.70 |
| 1,2-Dichloroethene, Total | ND | ug/ | 1 2.5 | 0.70 |
| Dibromomethane | ND | ug/ | J 5.0 | 1.0 |
| 1,2,3-Trichloropropane | ND | ug/ | 1 2.5 | 0.70 |
| Acrylonitrile | ND | ug/ | J 5.0 | 1.5 |
| Styrene | ND | ug/ | 1 2.5 | 0.70 |
| Dichlorodifluoromethane | ND | ug/ | 5.0 | 1.0 |
| Acetone | ND | ug/ | J 5.0 | 1.5 |
| Carbon disulfide | ND | ug/ | J 5.0 | 1.0 |
| 2-Butanone | ND | ug/ | J 5.0 | 1.9 |
| Vinyl acetate | ND | ug/ | J 5.0 | 1.0 |
| 4-Methyl-2-pentanone | ND | ug/ | J 5.0 | 1.0 |
| 2-Hexanone | ND | ug/ | J 5.0 | 1.0 |
| Bromochloromethane | ND | ug/ | 1 2.5 | 0.70 |
| 2,2-Dichloropropane | ND | ug/ | 1 2.5 | 0.70 |
| 1,2-Dibromoethane | ND | ug/ | 1 2.0 | 0.65 |
| 1,3-Dichloropropane | ND | ug/ | 1 2.5 | 0.70 |
| 1,1,1,2-Tetrachloroethane | ND | ug/ | l 2.5 | 0.70 |
| Bromobenzene | ND | ug/ | l 2.5 | 0.70 |
| n-Butylbenzene | ND | ug/ | l 2.5 | 0.70 |
| sec-Butylbenzene | ND | ug/ | l 2.5 | 0.70 |
| tert-Butylbenzene | ND | ug/ | 1 2.5 | 0.70 |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1810069

Report Date: 03/29/18

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/28/18 09:49

Analyst: PD

| Parameter | Result | Qualifier Units | s RL | MDL |
|-----------------------------|-----------------|-----------------|--------------|-------------|
| olatile Organics by GC/MS - | Westborough Lab | for sample(s): | 01-05 Batch: | WG1101340-5 |
| o-Chlorotoluene | ND | ug/l | 2.5 | 0.70 |
| p-Chlorotoluene | ND | ug/l | 2.5 | 0.70 |
| 1,2-Dibromo-3-chloropropane | ND | ug/l | 2.5 | 0.70 |
| Hexachlorobutadiene | ND | ug/l | 2.5 | 0.70 |
| Isopropylbenzene | ND | ug/l | 2.5 | 0.70 |
| p-Isopropyltoluene | ND | ug/l | 2.5 | 0.70 |
| Naphthalene | ND | ug/l | 2.5 | 0.70 |
| n-Propylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,3-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trichlorobenzene | ND | ug/l | 2.5 | 0.70 |
| 1,3,5-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,2,4-Trimethylbenzene | ND | ug/l | 2.5 | 0.70 |
| 1,4-Dioxane | ND | ug/l | 250 | 61. |
| p-Diethylbenzene | ND | ug/l | 2.0 | 0.70 |
| p-Ethyltoluene | ND | ug/l | 2.0 | 0.70 |
| 1,2,4,5-Tetramethylbenzene | ND | ug/l | 2.0 | 0.54 |
| Ethyl ether | ND | ug/l | 2.5 | 0.70 |
| trans-1,4-Dichloro-2-butene | ND | ug/l | 2.5 | 0.70 |
| | | | | |

Tentatively Identified Compounds

No Tentatively Identified Compounds

ND

ug/l



L1810069

Project Name: 123 HOPE STREET

Project Number: Report Date: 170357801 03/29/18

Lab Number:

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 03/28/18 09:49

Analyst: PD

| Parameter | Result | Qualifier | Units | RL | MDL | |
|-----------------------------------|------------|-------------|-------------|--------|-------------|--|
| Volatile Organics by GC/MS - West | borough La | b for sampl | e(s): 01-05 | Batch: | WG1101340-5 | |

| | Acceptance | | | | | | | |
|-----------------------|--------------|-------------------|--|--|--|--|--|--|
| Surrogate | %Recovery Qu | ualifier Criteria | | | | | | |
| 1,2-Dichloroethane-d4 | 96 | 70-130 | | | | | | |
| Toluene-d8 | 106 | 70-130 | | | | | | |
| 4-Bromofluorobenzene | 116 | 70-130 | | | | | | |
| Dibromofluoromethane | 89 | 70-130 | | | | | | |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1810069

Report Date: 03/29/18

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | %Recovery Qual Limits | RPD | RPD Qual Limits |
|--|------------------|------------|-------------------|--------------------------|-----|--------------------|
| Volatile Organics by GC/MS - Westborough | Lab Associated | sample(s): | 01-05 Batch: W | G1101340-3 WG1101340-4 | | |
| Methylene chloride | 81 | | 82 | 70-130 | 1 | 20 |
| 1,1-Dichloroethane | 89 | | 90 | 70-130 | 1 | 20 |
| Chloroform | 84 | | 84 | 70-130 | 0 | 20 |
| Carbon tetrachloride | 99 | | 99 | 63-132 | 0 | 20 |
| 1,2-Dichloropropane | 85 | | 84 | 70-130 | 1 | 20 |
| Dibromochloromethane | 82 | | 82 | 63-130 | 0 | 20 |
| 1,1,2-Trichloroethane | 84 | | 84 | 70-130 | 0 | 20 |
| Tetrachloroethene | 90 | | 88 | 70-130 | 2 | 20 |
| Chlorobenzene | 87 | | 85 | 75-130 | 2 | 20 |
| Trichlorofluoromethane | 100 | | 100 | 62-150 | 0 | 20 |
| 1,2-Dichloroethane | 82 | | 82 | 70-130 | 0 | 20 |
| 1,1,1-Trichloroethane | 96 | | 94 | 67-130 | 2 | 20 |
| Bromodichloromethane | 81 | | 83 | 67-130 | 2 | 20 |
| trans-1,3-Dichloropropene | 94 | | 93 | 70-130 | 1 | 20 |
| cis-1,3-Dichloropropene | 84 | | 85 | 70-130 | 1 | 20 |
| 1,1-Dichloropropene | 97 | | 96 | 70-130 | 1 | 20 |
| Bromoform | 79 | | 79 | 54-136 | 0 | 20 |
| 1,1,2,2-Tetrachloroethane | 86 | | 90 | 67-130 | 5 | 20 |
| Benzene | 87 | | 87 | 70-130 | 0 | 20 |
| Toluene | 90 | | 90 | 70-130 | 0 | 20 |
| Ethylbenzene | 95 | | 94 | 70-130 | 1 | 20 |
| Chloromethane | 94 | | 94 | 64-130 | 0 | 20 |
| Bromomethane | 51 | | 51 | 39-139 | 0 | 20 |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1810069

Report Date: 03/29/18

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | RPD Qual Limits | |
|--------------------------------|---------------------------|----------------|-------------------|------------|---------------------|-----|--------------------|--|
| Volatile Organics by GC/MS - W | estborough Lab Associated | sample(s): 01- | -05 Batch: W | G1101340-3 | WG1101340-4 | | | |
| Vinyl chloride | 98 | | 97 | | 55-140 | 1 | 20 | |
| Chloroethane | 100 | | 100 | | 55-138 | 0 | 20 | |
| 1,1-Dichloroethene | 91 | | 93 | | 61-145 | 2 | 20 | |
| trans-1,2-Dichloroethene | 89 | | 88 | | 70-130 | 1 | 20 | |
| Trichloroethene | 83 | | 84 | | 70-130 | 1 | 20 | |
| 1,2-Dichlorobenzene | 83 | | 82 | | 70-130 | 1 | 20 | |
| 1,3-Dichlorobenzene | 87 | | 87 | | 70-130 | 0 | 20 | |
| 1,4-Dichlorobenzene | 85 | | 86 | | 70-130 | 1 | 20 | |
| Methyl tert butyl ether | 85 | | 87 | | 63-130 | 2 | 20 | |
| p/m-Xylene | 90 | | 90 | | 70-130 | 0 | 20 | |
| o-Xylene | 90 | | 90 | | 70-130 | 0 | 20 | |
| cis-1,2-Dichloroethene | 83 | | 82 | | 70-130 | 1 | 20 | |
| Dibromomethane | 76 | | 79 | | 70-130 | 4 | 20 | |
| 1,2,3-Trichloropropane | 84 | | 92 | | 64-130 | 9 | 20 | |
| Acrylonitrile | 82 | | 82 | | 70-130 | 0 | 20 | |
| Styrene | 85 | | 85 | | 70-130 | 0 | 20 | |
| Dichlorodifluoromethane | 130 | | 120 | | 36-147 | 8 | 20 | |
| Acetone | 100 | | 110 | | 58-148 | 10 | 20 | |
| Carbon disulfide | 92 | | 92 | | 51-130 | 0 | 20 | |
| 2-Butanone | 91 | | 86 | | 63-138 | 6 | 20 | |
| Vinyl acetate | 83 | | 83 | | 70-130 | 0 | 20 | |
| 4-Methyl-2-pentanone | 83 | | 84 | | 59-130 | 1 | 20 | |
| 2-Hexanone | 92 | | 92 | | 57-130 | 0 | 20 | |
| | | | | | | | | |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number: L1810069

Report Date: 03/29/18

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits | |
|--------------------------------|---------------------------|------------|-------------------|-------------|---------------------|-----|------|---------------|--|
| Volatile Organics by GC/MS - W | estborough Lab Associated | sample(s): | 01-05 Batch: | WG1101340-3 | WG1101340-4 | | | | |
| Bromochloromethane | 77 | | 73 | | 70-130 | 5 | | 20 | |
| 2,2-Dichloropropane | 100 | | 100 | | 63-133 | 0 | | 20 | |
| 1,2-Dibromoethane | 83 | | 83 | | 70-130 | 0 | | 20 | |
| 1,3-Dichloropropane | 85 | | 86 | | 70-130 | 1 | | 20 | |
| 1,1,1,2-Tetrachloroethane | 89 | | 87 | | 64-130 | 2 | | 20 | |
| Bromobenzene | 83 | | 84 | | 70-130 | 1 | | 20 | |
| n-Butylbenzene | 120 | | 120 | | 53-136 | 0 | | 20 | |
| sec-Butylbenzene | 110 | | 110 | | 70-130 | 0 | | 20 | |
| tert-Butylbenzene | 88 | | 89 | | 70-130 | 1 | | 20 | |
| o-Chlorotoluene | 93 | | 94 | | 70-130 | 1 | | 20 | |
| p-Chlorotoluene | 95 | | 96 | | 70-130 | 1 | | 20 | |
| 1,2-Dibromo-3-chloropropane | 77 | | 81 | | 41-144 | 5 | | 20 | |
| Hexachlorobutadiene | 120 | | 110 | | 63-130 | 9 | | 20 | |
| Isopropylbenzene | 100 | | 100 | | 70-130 | 0 | | 20 | |
| p-Isopropyltoluene | 110 | | 110 | | 70-130 | 0 | | 20 | |
| Naphthalene | 90 | | 90 | | 70-130 | 0 | | 20 | |
| n-Propylbenzene | 100 | | 100 | | 69-130 | 0 | | 20 | |
| 1,2,3-Trichlorobenzene | 80 | | 81 | | 70-130 | 1 | | 20 | |
| 1,2,4-Trichlorobenzene | 83 | | 83 | | 70-130 | 0 | | 20 | |
| 1,3,5-Trimethylbenzene | 98 | | 98 | | 64-130 | 0 | | 20 | |
| 1,2,4-Trimethylbenzene | 100 | | 99 | | 70-130 | 1 | | 20 | |
| 1,4-Dioxane | 110 | | 140 | | 56-162 | 24 | Q | 20 | |
| p-Diethylbenzene | 100 | | 100 | | 70-130 | 0 | | 20 | |
| | | | | | | | | | |



Project Name: 123 HOPE STREET

Project Number: 170357801

Lab Number:

L1810069

Report Date:

03/29/18

| <u>Parameter</u> | LCS %Recovery | Qual | | .CSD ecovery | | %Recovery Limits | RPD | Qual | RPD Limits | |
|---|------------------|------------|-------|-----------------|-------------|---------------------|-----|------|---------------|--|
| Volatile Organics by GC/MS - Westborough La | ab Associated | sample(s): | 01-05 | Batch: | WG1101340-3 | WG1101340-4 | | | | |
| p-Ethyltoluene | 100 | | | 100 | | 70-130 | 0 | | 20 | |
| 1,2,4,5-Tetramethylbenzene | 96 | | | 96 | | 70-130 | 0 | | 20 | |
| Ethyl ether | 82 | | | 86 | | 59-134 | 5 | | 20 | |
| trans-1,4-Dichloro-2-butene | 90 | | | 90 | | 70-130 | 0 | | 20 | |

| | LCS | LCSD | Acceptance |
|-----------------------|----------------|----------------|------------|
| Surrogate | %Recovery Qual | %Recovery Qual | Criteria |
| 1,2-Dichloroethane-d4 | 98 | 98 | 70-130 |
| Toluene-d8 | 106 | 106 | 70-130 |
| 4-Bromofluorobenzene | 110 | 111 | 70-130 |
| Dibromofluoromethane | 91 | 91 | 70-130 |



Project Name: 123 HOPE STREET **Lab Number:** L1810069 Project Number: 170357801

YES

Report Date: 03/29/18

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Custody Seal Cooler

Α Absent

| Container Info | rmation | | Initial | Final | Temp | | | Frozen | |
|----------------|--------------------|--------|---------|-------|-------|------|--------|-----------|----------------|
| Container ID | Container Type | Cooler | рН | рН | deg C | Pres | Seal | Date/Time | Analysis(*) |
| L1810069-01A | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-01B | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-01C | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-02A | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-02B | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-02C | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-03A | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-03B | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-03C | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-04A | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-04B | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-04C | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-05A | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |
| L1810069-05B | Vial HCl preserved | Α | NA | | 3.1 | Υ | Absent | | NYTCL-8260(14) |



Project Name:123 HOPE STREETLab Number:L1810069Project Number:170357801Report Date:03/29/18

GLOSSARY

Acronyms

EDL - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated

values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis

of PAHs using Solid-Phase Microextraction (SPME).

EPA - Environmental Protection Agency.

LCS - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

LCSD - Laboratory Control Sample Duplicate: Refer to LCS.

LFB - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of

analytes or a material containing known and verified amounts of analytes.

MDL - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any

adjustments from dilutions, concentrations or moisture content, where applicable.

MS - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for

which an independent estimate of target analyte concentration is available.

MSD - Matrix Spike Sample Duplicate: Refer to MS.

NA - Not Applicable.

NC - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's

reporting unit.

NDPA/DPA - N-Nitrosodiphenylamine/Diphenylamine.

NI - Not Ignitable.

NP - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.

RL - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL

includes any adjustments from dilutions, concentrations or moisture content, where applicable.

RPD - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less

than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the

values; although the RPD value will be provided in the report.

SRM - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the

associated field samples.

STLP - Semi-dynamic Tank Leaching Procedure per EPA Method 1315.

TIC - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound

list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Footnotes

- The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

Data Qualifiers

A - Spectra identified as "Aldol Condensation Product".

- The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related

Report Format: DU Report with 'J' Qualifiers



В

Project Name:123 HOPE STREETLab Number:L1810069Project Number:170357801Report Date:03/29/18

Data Qualifiers

projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).

- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations
 of the analyte.
- E Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- G The concentration may be biased high due to matrix interferences (i.e, co-elution) with non-target compound(s). The result should be considered estimated.
- H The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I The lower value for the two columns has been reported due to obvious interference.
- M Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- NJ Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P The RPD between the results for the two columns exceeds the method-specified criteria.
- Q The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)
- **R** Analytical results are from sample re-analysis.
- **RE** Analytical results are from sample re-extraction.
- S Analytical results are from modified screening analysis.
- Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- ND Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.

Report Format: DU Report with 'J' Qualifiers



 Project Name:
 123 HOPE STREET
 Lab Number:
 L1810069

 Project Number:
 170357801
 Report Date:
 03/29/18

REFERENCES

Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - IV, 2007.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Alpha Analytical, Inc.
Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:**17873**

Revision 11 Published Date: 1/8/2018 4:15:49 PM

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Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624: m/p-xylene, o-xylene

EPA 8260C: <u>NPW</u>: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; <u>SCM</u>: lodomethane (methyl iodide), Methyl methacrylate, 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.

EPA 8270D: NPW: Dimethylnaphthalene,1,4-Diphenylhydrazine; SCM: Dimethylnaphthalene,1,4-Diphenylhydrazine.

EPA 300: <u>DW:</u> Bromide EPA 6860: <u>SCM:</u> Perchlorate

EPA 9010: NPW and SCM: Amenable Cyanide Distillation

SM4500: NPW: Amenable Cyanide, Dissolved Oxygen; SCM: Total Phosphorus, TKN, NO2, NO3.

Mansfield Facility

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87,101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

Drinking Water

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; EPA 353.2: Nitrate-N, Nitrite-N; SM4500NO3-F: Nitrate-N, Nitrite-N; SM4500F-C, SM4500CN-CE, EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B

EPA 332: Perchlorate; EPA 524.2: THMs and VOCs; EPA 504.1: EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT,SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, LACHAT 10-107-06-1-B: Ammonia-N, EPA 351.1, SM4500NO3-F, EPA 353.2: Nitrate-N, EPA 351.1, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D.

EPA 624: Volatile Halocarbons & Aromatics,

EPA 608: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan II, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625: SVOC (Acid/Base/Neutral Extractables), EPA 600/4-81-045: PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, SM9222D.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Be, Cd, Cr, Cu, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn.

EPA 245.1 Hg.

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

Document Type: Form Pre-Qualtrax Document ID: 08-113

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| E = NaOH F = MeOH $G = NaHSO_4$ $H = Na_2S_2O_3$ K/E = Zn Ac/NaOH O = Other Form No: 01-25 HC (rev. 3 | C = Cube O = Other E = Encore D = BOD Bottle | Relinquished | By: /44/ | 3/2' 3/25/18 | 71me 3/18 21/18 | Reg | Receive | d Byy DAA DZ3 | 31 | 133/ | Time (6) | start until any ambiguities are resolved. BY EXECUTING THIS COC, THE CLIENT HAS READ AND AGREES TO BE BOUND BY ALPHA'S TERMS & CONDITIONS. (See reverse side.) |