REMEDIAL INVESTIGATION REPORT FOR THE FORMER LUDLOW STREET WORKS SITE

Yonkers, New York

Site Number V00562

Prepared For:



Consolidated Edison Company of New York, Inc.

31-01 20th Avenue Long Island City, NY 11105

Prepared By:

PARSONS

Somerset, New Jersey

AUGUST 2014

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AUGUST 2014

"I ERIC GAULIN certify that I am currently a Qualified Environmental professional as defined in 6 NYCRR Part 375 and that this Remedial Investigation Report was prepared in accordance with all applicable statutes and regulations and in substantial conformance with the DER Technical Guidance for Site Investigation and Remediation (DER-10)."

Signature

Date

TABLE OF CONTENTS

	<u>P</u>	age
SECTION 1	INTRODUCTION	1
1.1	Report Organization	1
SECTION 2	SITE BACKGROUND	2
2.1	Site Overview	2
2.2	Adjoining Property Description	2
2.3	Site History	3
2.4	Topography, Regional Geology, and Hydrogeology	3
2.5	Previous Investigations/Remedial Measures	4
SECTION 3	REMEDIAL INVESTIGATION ACTIVITIES	6
3.1	Site Inspection and Preliminary Investigation Activities	6
3.2	Utility Clearance	6
3.3	Test Pit Excavation	7
3.4	Soil Boring Installation	7
3.5	Monitoring Well Installation and Development	8
3.6	Surveying	9
3.7	Groundwater Sampling	.10
3.8	Soil Gas Investigation Activities	.10
3.9	Tidal Study	.11
3.10	Management of Investigation-Derived Waste	.12
3.11	Data Validation and Reporting	.12
SECTION 4	REMEDIAL INVESTIGATION RESULTS	.13

	4.1	Site Geology	13
	4.2	Site Hydrogeology	13
	4.3	Former MGP Structures	13
	4.4	Analytical Results AND fIELD oBSERVATIONS	15
	4.4.1	Subsurface Soil	15
	4.4.2	Groundwater	18
	4.4.3	Soil Gas Sample Results	22
SE	CTION 5	EXTENT OF MGP NAPL IMPACTS	24
	5.1	MGP NAPL – 6 to 11 feet	24
	5.2	MGP NAPL – 11 to 21 Feet	24
	5.3	MGP NAPL – 21 to 25 Feet	25
	5.4	MGP NAPL – DEEPER THAN 35 38 Feet	25
SE	CTION 6	EXPOSURE ASSESSMENT	26
	6.1	SOILS – VOCs, SVOCs and MGP NAPL	26
	6.2	SOILS – INORGANIC IMPACTS	26
	6.3	GROUNDWATER – VOCs, SVOCs and MGP NAPL	27
	6.3	GROUNDWATER – Fill Related Constituents	27
	6.4	SOIL VAPOR	27
SE	CTION 7	CONCLUSIONS AND RECOMMENDATIONS	28
	7.1	Conclusions	28
	7.2	Recommendations	30
SE.	CTION 8	REFERENCES	31

LIST OF TABLES

- Table 1 RI Sample Summary
- Table 2 Summary of Groundwater Elevations
- Table 3 Summary of Soil Analytical Data
- Table 4 Summary of Groundwater Analytical Data
- Table 5 Summary of Soil Gas Analytical Data

LIST OF FIGURES

Figure 1	Site Location Map	
Figure 2	Former MGP Structures	
Figure 3	Sample Location Map	
Figure 4	Summary of VOC Exceedances in Soil	
Figure 5	Summary of SVOC Exceedances in Soil	
Figure 6	Summary of Inorganic Exceedances in Soil	
Figure 7A	Cross Sections A-A', B-B'	
Figure 7B	Cross Sections C-C', D-D'	
Figure 7C	Cross Section E-E'	
Figure 8	Groundwater Elevation Contour Map	
Figure 9	Summary of VOC Exceedances in Groundwater	
Figure 10	Summary of SVOC Exceedances in Groundwater	
Figure 11	Summary of Inorganic Exceedances in Groundwater	
Figure 12	Approximate Location of MGP NAPL – 6 to 11 Feet	
Figure 13	Approximate Location of MGP NAPL – 11 to 21 Feet	
Figure 14	Approximate Location of MGP NAPL – 21 to 25 Feet	
Figure 15	Approximate Location of MGP NAPL – 36 to 38 Feet	

LIST OF APPENDICES

Appendix	Α	Test Pit	[_00
1 ippelial/i	1 L	I OSC I IC	

Appendix B Soil Boring and Monitoring Well Logs

Appendix C Groundwater Sampling Logs

Appendix D Tidal Study Results

Appendix E Soil Gas Laboratory Analytical Report

Appendix F Data Usability Summary Reports

INTRODUCTION

This Remedial Investigation Report (RIR) presents the results of the remedial investigation (RI) of the former Ludlow Street Works former manufactured gas plant and holder station (MGP) located in Yonkers, New York. The RI was conducted between October 2012 and April 2013 by Parsons on behalf of Consolidated Edison Company of New York, Inc. (Con Edison). The goal of the RI was to determine the extent of impacts related to former MGP operations identified during the previous Site Characterization Study and summarized in the *Site Characterization Report* (Parsons, 2011). RI activities were completed in accordance with the New York State Department of Environmental Conservation's (NYSDEC) approved *Remedial Investigation Work Plan* (Parsons, 2011) and consistent with processes outlined in the NYSDEC Technical Guidance for Site Investigation and Remediation (NYSDEC DER-10, 2010).

1.1 REPORT ORGANIZATION

The RI was conducted by Parsons. The field investigation activities and results are documented in the following sections:

- Section 1: Introduction
- Section 2: Site Background
- Section 3: Remedial Investigation Activities
- Section 4: Remedial Investigation Results
- Section 5: Extent of MGP NAPL Impacts
- Section 6: Exposure Assessment
- Section 7: Conclusions and Recommendations
- Section 8: References

SITE BACKGROUND

2.1 SITE OVERVIEW

Con Edison has entered into a Voluntary Cleanup Agreement with the NYSDEC to investigate, and if necessary, remediate potential impacts at former MGP properties. One of these facilities, the former Ludlow Street Works (Voluntary Cleanup Agreement Index No. D2-0003-02-08; Site ID No. V00562), a former MGP and holder station, was operated between 1879 and 1936 by predecessor companies of Con Edison.

The grounds of the former Ludlow Street MGP Works Site (Site) occupy present-day tax Block 171 Lot 17, which is owned by the City of Yonkers. This approximately 1.35-acre parcel is a roughly rectangular plot with the long axis trending north-south and is bound on the north by Downing Street, on the south by the western-end of Knowles Street and along the west side by a Metro North Railroad right-of-way (ROW). Two separate tax lots that are owned by others form the eastern Site boundary. These parcels include a vacant lot (Block 171 Lot 11) along the northern half of the eastern boundary. This parcel is owned by the City of Yonkers, but there are no apparent municipal operations there. The parcel along the southern half of the eastern boundary is used as a commercial metal fabrication shop.

Currently, there is only one small single-story concrete block structure located on the southeastern corner of the southern portion of the Site. The northern and southern portions of the Site are separated by a concrete block retaining wall. The southern portion of the Site, where the majority of the former MGP operations took place and where two of the former holder structures were located, is currently used by the City of Yonkers Department of Public Works (DPW) as a bulk storage and staging area for road salt and associated equipment (e.g., trucks, front-end loaders, etc.) (Figure 1). The surface of the active DPW yard consists of gravel and dirt. Although the surface of DPW yard was formerly covered with bituminous asphalt paving this surface is now generally absent, likely due to its inadvertent removal during salt load-out operations. The southeast corner of northern portion of the site contained the 50,000 cu ft holder and a portion of the former Retort house was located in its southwest area. The elevation of the northern or upper portion of the Site is approximately 10 to 26 feet higher than the southern or lower portion (Chazen, 2013) as a result of City of Yonkers' placement there of various debris. The City also removed trees and brush from the elevated northern portion of the site and subsequently leveled the surface there.

The Site is located within a mixed-use industrial district of Yonkers that also contains residential and commercial-use properties. The Site fronts two streets, at approximately 162 Downing Street and the intersection of Knowles and Bridge Streets in Yonkers, New York. The Site consists of approximately 1.35 acres and is identified as Section 1, Block 171, Lot 1 (Figure 2).

2.2 ADJOINING PROPERTY DESCRIPTION

The Site is bordered to the north by Downing Street; to the east by a metal fabrication shop (the Waverly property); to the south by Knowles Street (formerly Fernbrook Street) and a Dunkin Donuts bakery; and to the west by the Metropolitan Transportation Authority

(MTA)/New York Central Railroad railway. The Site is located approximately 700 feet east of the Hudson River (*GEI*, 2003). However, based on 1886 Sanborn Fire Insurance map for the area where the Site was located, the shore of the Hudson River when the former MGP first began operation was located along the west side of the railroad tracks. A ccordingly, at that time, the Site was located less than 100 feet from the Hudson River shoreline.

2.3 SITE HISTORY

Historical research of the Site was previously conducted and is documented in the *Manufactured Gas Plant History, Ludlow Street Works, Yonkers, New York (GEI, 2003)* which indicates that the Site was operated as an MGP and gas holder station by several predecessor companies of Con Edison on the southern portion of the Site between 1879 and 1936 and on the northern portion between 1923 and 1936. The Municipal Gas Company of Yonkers (i.e., Yonkers Municipal Gas Company) acquired the southern portion of the Site in 1879 from a private owner. The Westchester Lighting Company acquired the northern portion of the Site in 1923. The Municipal Gas Company constructed an MGP at the Site and began manufacturing gas around 1880.

In 1886, there were three buildings, two gas holders, and three underground naphtha storage tanks located on the southern portion of the Site (Figure 2). One holder was located on the southern portion of the Site and had a 22,800 cubic feet (cu ft) capacity. The other holder was located on the northern portion of the Site and had a 50,000 cubic cu ft capacity. One building, which contained the lime house, purifying house, and meter room, was located at the southwest corner side of the Site. To the north of this building was the engine room and retorts building. A boiler house was located at the southeastern corner of the Site. The underground naphtha storage tanks were located approximately at the middle of the west side of the Site. An underground pipe (shown on the 1886 Sanborn Fire Insurance Map) was used to pump naphtha from boats on the Hudson River to the tanks. By 1898, the plant was no longer manufacturing gas, and became a holder station for the storage of natural gas.

A third larger gas holder (250,000 cu ft) was constructed before 1907, replacing the building that contained the lime house, purifying house, and meter room and a portion of the retorts and engine room building. The two smaller holders were removed from the Site between 1917 and 1931. By 1936, the Site was no longer used as a holder station. The Site use between 1936 and 1946 is unknown (*GEI*, 2003). The Site was sold in 1946 and was owned by various bus companies for bus storage and parking until 1973 when the City of Yonkers acquired the Site. The City of Yonkers has used the Site as a DPW storage yard for vehicles and trucks, road salt, materials, street sweeping dumping, debris, fuel tanks, asbestos-containing materials, 55-gallon drums, and other unidentified waste materials since 1973.

2.4 TOPOGRAPHY, REGIONAL GEOLOGY, AND HYDROGEOLOGY

The Site is located in the southwestern corner of Yonkers, approximately 700 feet east of the Hudson River. B ased on the Site survey performed in December 2012, the ground surface elevation of the southern (or lower) portion of the Site is approximately 16 feet above mean sea level (MSL) and the elevation of the surveyed northern (or upper) portion of the Site is between approximately 38 and 44 feet above MSL. The Site is located in the Manhattan Prong section of the New England Uplands physiographic region of New York and is near a northeast-trending geologic contact between the Fordham Gneiss and the Inwood Marble. During the RI, bedrock

at the Site was encountered from approximately 39 to 50 feet below ground surface (feet bgs) which translates into approximately 3.4 feet below MSL at SB-14 and 23.2 feet below MSL at SB-12. The bedrock is overlain by glacial till deposits, alluvium, and fill material (*GEI*, 2003). Based on groundwater levels measured in monitoring wells during the RI between October 2012 and January 2013, the groundwater elevation at the Site ranged between 2.84 and 5.79 ft above MSL. Groundwater generally flows to the west-northwest, toward the Hudson River. As indicated in earlier reports and confirmed by the tidal study conducted during this RI, groundwater levels are minimally influenced by tidal fluctuation of water in the adjoining Hudson River.

2.5 PREVIOUS INVESTIGATIONS/REMEDIAL MEASURES

Two investigations were previously conducted on the Site. The first was completed between 1993 and 1997 and the second in 2010. These investigations are summarized below.

In 1993, Vollmuth & Brush of Blue Point, New York performed a two-phase sampling and analysis program on the Site and the adjacent lots (Vollmuth and Brush, 1997). A report entitled "Phase I and II Sampling and Analysis Program 151-154 Downing Street, DPW Garage, City of Yonkers, Yonkers, NY" was prepared for the Department of Planning and Development of City of Yonkers by Vollmuth & Brush in 1997 (Vollmuth and Brush, 1997). The first phase of the investigation entailed a soil gas survey, and the second phase entailed the sampling and analysis of subsurface soils. Eleven soil samples were collected from inside the former DPW garage building on the eastern abutting property and on the Site. Soil samples were analyzed for volatile organic compounds (VOCs) and semivolatile organic compounds (SVOCs) from depths of 0 to 2 or 2 to 4 feet below grade. Soils samples were analyzed for metals from depths of 0 to 2, 2 to 4, 6 to 8, or 10 to 12 feet below grade. Vollmuth & Brush compared the detected concentrations of the VOCs and SVOCs to their corresponding NYSDEC Soil Cleanup Objectives (SCOs). Based on these comparisons, it was determined that concentrations of all VOCs and SVOCs detected were below their respective SCOs. Five metals, namely copper, iron, lead, selenium, and zinc, were detected at concentrations above their respective SCOs in at least one soil sample.

In 2010, Con Edison contracted Parsons to conduct a Site Characterization Study (SCS) of the site. The scope of Site Characterization activities included the excavation of five test pits (designated TP-01 through TP-05 and six soil borings (designated SB-1, SB-2, SB-4 through SB-7), and installation of five monitoring wells (designated MW-1 through MW-5). Since historical information indicated that operations of the former MGP were primarily located at the southern (or lower) portion of the Site, field investigation activities focused on these areas of the former site and the adjacent sidewalk. During the SCS, remnants of several former MGP structures were identified on the property. Where present, coal tar was generally detected as residue or at saturated levels as Non-Aqueous Phase Liquid (NAPL) in the vicinity of the former/remnant structures. Fingerprint analysis of coal tar NAPL samples verified that the tar residues were formed as a result of the production of manufactured coal gas. Contaminants potentially related to former MGP activities were also detected in soil samples collected during the SCS at concentrations exceeding their respective NYSDEC SCOs. These contaminants were generally encountered at or below the groundwater table (e.g., between 10.3 ft to 16.1 feet bgs) and extended to depths ranging from 30 ft to 32 feet bgs. The vertical extent of these contaminants was delineated at all sampling locations, with the exception of SB-02.

Contaminants potentially related to former MGP activities including VOCs, SVOCs and metals were also detected in groundwater at concentrations exceeding the NYSDOH Ambient Water Quality Standards, Criteria and Guidance Values (AWQSGV) for Class GA water.

Additionally, elevated concentrations of sodium, manganese, and magnesium were detected in Site soil and groundwater and are suspected to be associated with use of the Site as a road salt storage facility, or due to the presence of urban fill materials.

Based on the findings of the SCS, it was recommended by Parsons in the SCS Report (*March 2011*) that further investigation of the site be conducted to delineate the impacts identified. Subsequently, a Remedial Investigation Work Plan (Parsons, 2011) was then developed and approved by the NYSDEC.

REMEDIAL INVESTIGATION ACTIVITIES

The following sections describe the field activities conducted by Parsons between October 2012 and April 2013 in accordance with the NYSDEC approved *Remedial Investigation Work Plan (Parsons, 2011)*. The RI field investigation included a geophysical survey, utility clearance activities, excavation of a test pit, advancement of soil borings, installation and development of groundwater monitoring wells, sampling and laboratory analysis of soil, groundwater and soil vapor samples, tidal study, and surveying. During all intrusive activities, a Community Air Monitoring Plan (CAMP) was implemented in accordance with the approved work plan and associated site-specific Health and Safety Plan (HASP). Sampling locations are depicted on Figure 3. Table 1 provides a summary of the samples and analyses.

3.1 SITE INSPECTION AND PRELIMINARY INVESTIGATION ACTIVITIES

On October 15, 2012, a Site inspection was conducted with Con Edison, Parsons, representatives of the City of Yonkers DPW and Westchester County Sewer Department to review the proposed sampling locations (borings, wells, test pits, etc.). Proposed sampling locations were adjusted during the site inspection, as necessary, based on Site conditions, access, utilities, and or safety considerations. All changes were made in consultation with Con Edison and the NYSDEC. Specific modifications to sampling locations are further discussed below.

3.2 UTILITY CLEARANCE

The utility clearance was performed as follows:

- A Code 753 mark out was requested by Aquifer Drilling & Testing, Inc. (ADT) of Mineola, New York;
- Site reconnaissance was performed, followed by meetings with the Yonkers Department of Public Works, the Westchester County Department of Environmental Facilities, and the owners of the Waverly property;
- Utility maps were reconciled with the mark outs, and with the representatives from these parties;
- A subsurface geophysical survey was performed by Diversified Geophysics, Inc. (DGI), of Mineola, New York at all proposed boring, monitoring well and test pit locations;
- The 54-inch forced sewer main on the MTA parcel was exposed prior to boring advancement;
- Utility clearance key-hole test pits were excavated to 5 feet at each proposed soil boring, monitoring well, and soil vapor location for subsurface utilities.

During these excavation activities, soils were screened for VOCs using a photoionization detector (PID), their physical characteristics (e.g., soil type, grain size, color, etc.) were described, and notes of any evidence of physical impacts observed (staining, odor, sheen, NAPL, etc.) were recorded.

3.3 TEST PIT EXCAVATION

One investigation test pit (TP-6) was excavated in an effort to determine if the former 4-inch naphtha pipe shown on the 1886 S anborn Fire Insurance map was still present, and, if so, determine if there are any potential impacts from this pipe to the surrounding soil and or groundwater. Test pit excavation was completed on November 7, 2012. Test pit excavation was conducted by ADT under the supervision of a Parsons geologist. The test pit location is depicted on Figure 3. Based on Site conditions and observations made during implementation of the RI, TP-6 was excavated 10 feet west of its originally proposed location due to the presence of a 54-inch diameter sanitary sewer force main. The test pit was advanced with a rubber-tired backhoe to six feet bgs, at which point, groundwater was encountered.

During test pit excavation, the excavated soil and fill were temporarily placed on impervious plastic sheeting adjacent to the test pit. The test pit excavation was conducted in 6-inch lifts to its final depth. The excavated material showed no apparent evidence of impacts and was placed back into the excavation in reverse order from which it was removed. The lower portion of the test pit was backfilled in 12-inch lifts. The upper portion of the test pit was backfilled in 6-inch lifts using suitable excavated materials (e.g. material that did not contain deleterious materials, saturated or near saturated fill, rocks, soils clumps bigger than 6 inches in size, etc.).

While the test pit remained open, subsurface conditions were logged and photographed, the test pit walls and floors were inspected for evidence of MGP-related impacts (e.g., odors, staining, sheens, NAPL, and or PID readings above background levels) and the dimensions of the test pit and any subsurface features were measured. All field observations and measurements are summarized in the test pit log provided in Appendix A.

The soil sample, designated TP-6 Floor(x-y) was collected from the test pit was submitted to Chemtech Consulting Group, Inc of Mountainside, NJ (Chemtech) (NYS DOH, ELAP Certified) for analysis of Target Compound List (TCL) VOCs by EPA Method 8260, T CL SVOCs by EPA Method 8270, TAL metals by EPA Method 6000/7000 Series, and cyanide by EPA Method 9012. The analytical results for this sample are summarized in Table 3.

3.4 SOIL BORING INSTALLATION

A total of eight soil borings (SB-9 through SB-14, SB-16, and SB-17) were advanced during the RI to further delineate the nature and extent of MGP impacts in subsurface soils. The soil borings were completed between October 25, 2012 and December 21, 2012. Soil borings were advanced by ADT under the supervision of a Parsons geologist. Soil borings were completed to depths ranging from approximately 25 to 53 feet bgs, depending on observed impacts and or refusals. Soil boring locations are shown on Figure 3 and the corresponding boring logs are presented in Appendix B. Based on Site conditions and observations made during implementation of the RI activities, the following additions/modifications were made to the work plan:

- SB-9/SV-1 was relocated approximately 10 feet east due to the presence of a 30-inch sewer line in the vicinity of its original location;
- SB-11 was relocated approximately 10 feet east due to the presence of a 54-inch forced sewer main in the vicinity of its original location;

- SB-12 was relocated approximately 5 feet south due to its proximity to an unstable debris pile slope;
- SB-13 was added to the scope in order to determine the presence, nature, and extent of NAPL impacts encountered at SB-10 from 27 to 42.5 feet bgs;
- SB-14 was added to the scope in order to determine the presence, nature, and extent of NAPL impacts encountered at SB-13 from 47 to 49 feet bgs;
- SB-16 was added to the scope in order to determine the presence, nature, and extent of NAPL impacts encountered at SB-10 and SB-12; and
- SB-17 was added to the scope in order to determine the presence, nature, and extent of NAPL impacts encountered at MW-7 from 35 to 40 and SB-13 from 47 to 49 feet bgs.

Soil borings were advanced using a track-mounted sonic drilling rig. Soil core was collected continuously to the bottom of the boring using 5-foot and 10-foot long, 4-inch diameter stainless steel soil core samplers. The soil core was screened for the presence of VOCs using a PID and logged. Physical characteristics of each sample were recorded (e.g., soil type, color, texture, moisture content, etc.), along with physical evidence of any impacted material (e.g., oil-like or tar-like NAPL, staining, sheens, odors, etc).

Subsequent to characterizing the core, soil samples were collected from discrete intervals based on the criteria listed below. The samples were placed in clean sample jars which were stored in a cooler containing ice prior to shipment to Chemtech for analysis.

- One sample was collected from the zone with the highest PID readings or strong evidence of other physical impacts (e.g., odor, staining, sheen, etc.). If physical impacts or elevated PID readings were not detected, then a sample was collected from the upper portion of the boring or directly above the water table (if present).
- One sample was collected below the impacted zone (if present) or near the base of the boring to identify the vertical extent of any impacts at the location.

The soil samples were submitted to Chemtech of Mountainside, NJ and analyzed for TCL VOCs, TCL SVOCs, cyanide, and TAL metals. A summary of the soil samples collected and analyses performed is provided in Table 1.

Upon completion, the soil borings were grouted with cement-bentonite grout using a tremie pipe. Drilling equipment was decontaminated between each boring. D rill cuttings and decontamination water were containerized in new USDOT approved 55-gallon steel drums and handled as described in Section 3.10

3.5 MONITORING WELL INSTALLATION AND DEVELOPMENT

A total of three monitoring wells (MW-7, MW-8, and MW-9) were installed during the RI activities to further delineate the nature and extent of MGP impacts in groundwater. Monitoring wells were installed from October 26, 2012 through December 19, 2012. Monitoring wells were advanced by ADT under the supervision of a Parsons geologist. Monitoring wells were advanced to depths ranging from approximately 25 to 53 feet bgs, depending on observed impacts and refusals; monitoring wells were constructed at depths ranging from 14 to 50 feet bgs. Monitoring well locations are shown on Figure 3 and the corresponding boring/well logs

are presented in Appendix B. B ased on Site conditions and observations made during implementation of the RI activities, the following additions/modifications were made to the work plan:

- MW-2 was added to the scope in order to provide a replacement of the original MW-2 which was believed to be destroyed by DPW site activities;
- MW-5 was added to the scope in order to provide a replacement of the original MW-5 which was believe to be destroyed by DPW site activities;
- MW-6/SV-2 was not installed due to the presence of a 30-inch sewer line in the vicinity of its original location. Due to Site conditions and safety concerns, it was determined that there was no relocation option for this monitoring well/soil vapor location; and
- MW-9 was relocated approximately 18 feet southeast due to the presence of a 54-inch forced sewer main in the vicinity of its original location;

Monitoring well borings were completed as described above in Section 3.4. The monitoring well screens were installed at depths ranging from 4 to 50 feet bgs. Soil samples were not collected from the replacement wells MW-2 and MW-5.

The monitoring wells were constructed with 2-inch inner diameter, threaded, flush-joint, PVC casing and 10 to 20 feet lengths of 0.02-inch slot screen. The annular space around each well screen was backfilled with a No. 2 clean sand filter pack extending from the bottom of the well to at least 2 feet above the top of the screen. The annular space around the well riser was sealed with at least 2 feet of hydrated bentonite pellets on top of the sand pack. The remainder of the boring was backfilled with cement-bentonite grout to approximately 1 foot below grade. Monitoring wells MW-2, MW-5, MW-8 and MW-9 were finished with a locking, flush-mount box set in concrete. Monitoring well MW-7 was finished with a protective stick-up casing set in concrete.

Monitoring well development was conducted on November 9, 2012, D ecember 26, 2012, and January 3, 2012 a minimum of 24 hours after installation. Each new monitoring well was developed until reasonably free of sediment and until turbidity measurements were less than 50 nephelometric turbidity units (NTUs), if possible, and/or until the measurements of water quality parameters [i.e., pH, temperature, and conductivity] stabilized. The water quality parameters were measured and recorded approximately every 5 minutes until stabilization was achieved. Stabilization is considered to be achieved when three successive measurements of the water quality parameters were within 10%. It is noted that stabilization was not achieved at monitoring well MW-2 as it ran dry numerous times. Non-disposable drilling equipment was decontaminated between well locations. Monitoring well drill cuttings, well development water, and decontamination water were containerized in new 55-gallon steel drums and handled as described in Section 3.10.

3.6 SURVEYING

At the conclusion of drilling activities, Chazen, a licensed New York state land surveyor mobilized to the Site and surveyed the horizontal and vertical location of each new soil boring, monitoring well, and test pit location. Other pertinent planimetric features of the Site, such as manholes, building walls, etc. were also surveyed. Two elevation measurements were taken at each well location to identify the top of the PVC casing and the grade elevation, respectively.

The survey elevations were measured to an accuracy of 0.01 feet above the National Geodetic Vertical Datum of 1988 (NGVD 1988). Horizontal coordinates were surveyed relative to the New York State Plane North American Datum of 1983 (NAD83).

3.7 GROUNDWATER SAMPLING

On November 26 and 27, 2012, g roundwater samples were collected from six monitoring wells (MW-1, MW-3, MW-4, MW-7 through MW-9). Monitoring wells MW-2 and MW-5 were sampled during the second round of groundwater sampling based on their date of installation.

On April 5 and 8, 2013, groundwater samples were collected from seven monitoring wells (MW-1 through MW-5, MW-7, and MW-9). Monitoring well MW-8 was not sampled during the second round of groundwater sampling due to access issues.

Groundwater sampling was conducted in accordance with the January 19, 2010 USEPA Region 1 Low Stress (Low Flow) Purging and Sampling Procedure for the Collection of Groundwater Samples from Monitoring Wells. P rior to collecting samples, the depths to groundwater, top of coal tar NAPL (if any) and bottom of well were measured relative to the top of well casing using an electronic oil/water interface probe accurate to 0.01 f oot. Table 2 provides a summary of groundwater level measurements and elevations taken on November 19, 2012, January 3 and April 5, 2013.

Prior to purging and immediately after removing the locking well cap, the concentration of total VOCs concentrations in the headspace air of each well was measured using a PID. Each well was purged using a submersible pump and low-flow purging techniques with dedicated tubing until stabilization of water quality parameters (including temperature, conductivity, pH, dissolved oxygen, ORP, and turbidity was achieved, and thus allow for the collection of a representative groundwater sample. As discussed in Section 3.7, monitoring well MW-4 ran dry a number of times prior to the stabilization of water quality parameters. During the purging process, water quality parameters were measured approximately every 5 minutes.

Water quality parameter measurements and observations were recorded on groundwater sampling field forms, which are provided in Appendix C. Groundwater samples were submitted to Chemtech Laboratories for analysis of TCL VOCs, TCL SVOCs, TAL Metals and total cyanide. During the April 2013 sampling event, groundwater samples MW-1 and MW-4 were filtered and analyzed for dissolved TAL metals, due to the high turbidity observed in groundwater from these wells. Non-dedicated sampling equipment (e.g., oil/water interface probe, submersible pump) was decontaminated between wells. Water generated during well purging and equipment decontamination was placed in 55-gallon drums and handled as described in Section 3.10.

3.8 SOIL GAS INVESTIGATION ACTIVITIES

To evaluate potential presence of VOCs in soil gas at the site and adjacent parcels, soil gas samples were collected from two locations (designated SV-3 and SV-4) adjacent to the one-story building bordering the southern boundary of the former site and one location, (designated SV-1) adjacent to the one story metal building located on the adjoining property along the eastern Site boundary. The soil gas sampling locations are shown on Figure 3. Soil gas samples were collected in accordance with the New York State Department of Health (NYSDOH) Final Guidance for Evaluating Soil Vapor Intrusion (NYSDOH, 2006).

Prior to soil gas sampling, the approximate depth of the bottom of the basement adjacent to the SV-3 and SV-4 locations (Dunkin Donuts Bakery), and the approximate depth of the building foundation slab adjacent to the SV-1 (Waverly Property) locations relative to the ambient ground surface were determined. The slab bottom was estimated to be 6 feet at the eastern building and the basement was estimated to be at 7 feet bgs at the southern building. Soil gas samples were collected from one foot below the bottom of each slab or basement bottom. A ccordingly, the sample from the eastern (Waverly) property (i.e., SV-1) was collected from a depth of seven feet bgs and the samples from the southern (Dunkin Donuts Bakery) property (i.e., SV-3 and SV-4) were collected from a depth of eight feet bgs.

Soil gas samples were collected using a probe/sampling assembly comprised of stainless steel tubing fitted with an expendable steel probe point. Prior to soil gas sampling, the sampling probe was driven into the sub-slab soil via a direct push method to the desired sampling depth. The expendable point was released into the soil and the sampler was withdrawn approximately 0.5 feet to create a void space. The space between the probe rod and the concrete slab was filled with granular bentonite, which was then hydrated to form an airtight seal to prevent aboveground ambient air from influencing the quality of a subsurface soil gas sample. A minimum of one implant volume (the volume of the sample probe and tube) was purged prior to sample collection.

A tracer gas (helium) was utilized in the field to verify the integrity and effectiveness of the hydrated-bentonite seal prior to sample collection at select locations. An enclosure (inverted plastic pail) was placed on the concrete slab over the location where the probe intersected the ground and the atmosphere within the enclosure was enriched with helium gas. A fter the atmosphere within the enclosure had been enriched, a vapor sample from the probe was collected and screened in the field with a helium detector to ensure the tracer gas was absent (reading of 0 PPM). The soil gas samples were collected into laboratory-prepared 6-liter Summa canisters equipped with calibrated flow regulators and vacuum gauge. The vacuum gauge was used to measure the initial and final vacuum in the canisters. Flow rates for both purging and sample collection did not exceed 0.2 liters per minute. All soil gas samples were submitted to Air Toxics LTD of Folsom, CA for laboratory analysis of VOCs using modified EPA Method TO-15. A summary of the soil gas samples collected and analyses performed is provided in Table 1. Soil gas laboratory results are provided in Appendix E.

3.9 TIDAL STUDY

A study was conducted at five monitoring wells to evaluate the potential influence of tidal fluctuations in the nearby Hudson River on groundwater elevations at and down gradient from the Site.

The tidal study involved the installation of Micro-DiverTM pressure transducer/data loggers in monitoring wells MW-1, MW-4, MW-7, MW-8, and MW-9. The transducers were set to record groundwater levels every 10 second intervals over a 72-hour period. The data was downloaded from the transducers and converted to elevation values. Graphs summarizing the groundwater elevation data were prepared and are provided in Appendix D.

3.10 MANAGEMENT OF INVESTIGATION-DERIVED WASTE

Investigation-derived waste (IDW), which included decontamination wash and rinse water, soil cuttings, purge water, debris, and used personal protective equipment (PPE), was containerized in Department of Transportation (DOT)-approved 55-gallon drums. The drums were sealed at the end of each work day and labeled with the date, the well or boring number(s), and the type of waste (e.g., drill cuttings, purge water). Parsons collected representative waste characterization samples of the IDW and coordinated transportation and disposal. Clean Earth of North Jersey, Inc. of Kearny, New Jersey disposed of the Site IDW at an offsite Con Edison-approved location in accordance with applicable local, state, and federal regulations.

3.11 DATA VALIDATION AND REPORTING

Data validation was performed in accordance with the USEPA Region II standard operating procedures (SOPs) for organic and inorganic data review which were in effect at the time of data validation (USEPA 2006; 2008a.2008b). These validation guidelines are regional modifications to the National Functional Guidelines for organic and inorganic data review (USEPA, 1999 and 2004). Validation included the following:

- Verification of 100% of all quality control (QC) sample results (both qualitative and quantitative);
- Verification of the identification of 100% of all sample results (both positive hits and non-detects);
- Recalculation of 10% of all investigative sample results; and
- Preparation of a Data Usability Summary Report (DUSR).

The quality of the data has been assessed and is documented in the DUSR provided in Appendix F. In summary, the results of the data usability assessment show that the collected analytical data for soil, groundwater and soil gas are valid for the intended purposes of the RI.

REMEDIAL INVESTIGATION RESULTS

This section presents the results of the RI. The goal of the RIis to determine the extent of impacts related to former MGP operations identified during the previous Site Characterization Study. Field observations are summarized followed by a review of the analytical data for the soil, groundwater, and soil gas samples.

4.1 SITE GEOLOGY

The geology encountered in the soil borings and test pits during the RI is summarized in the logs provided in Appendices A and B, and are interpreted in cross sections presented on Figures 7A, 7B, and 7C. The logs show that soil in the upper 5 feet (MW-8) to 12 feet (SB-11) west of the railway, upper 15 to 23 feet in the south and central portions and upper 26 to 38 feet in the north portions of the Site is characterized as urban fill. This material generally consists of sand, gravel and silt with various debris including cobbles, cement, brick, metal, wood fragments, etc. Natural deposits of fine to coarse-grained sand with some silt were encountered beneath the fill. These deposits were ranged from approximately 5 to 35 feet thick. Bedrock at the Site was encountered at depths ranging from approximately 39 to 50 feet bgs, which translates into approximately 3.4 feet below MSL at SB-14 (i.e., upper / northern portion of the site) and 23.2 feet below MSL at SB-12 (i.e., central portion of the site)

4.2 SITE HYDROGEOLOGY

A complete site-wide gauging event was conducted on January 3, 2013. The groundwater levels and corresponding elevations are summarized in Table 2. The January 3, 2013 groundwater elevations were used to produce a Site groundwater contour map (Figure 8). Groundwater contours based on the January 2013 gauging event indicate that the groundwater flow direction is predominantly to the west-northwest towards the Hudson River.

Additionally, the results of the tidal study indicate that groundwater elevations in monitoring wells MW-7, MW-8, and MW-9 are minimally influenced by tidal fluctuations in the Hudson River. The groundwater elevations in these wells showed coincident changes with the tides of less than 0.1 feet per tidal cycle. Such a small influence is not suspected to cause a change in groundwater flow directions in this area. With the exception of MW-7, response of groundwater elevations to tidal fluctuations were not detected in any of the on-site wells. Tidal study data and graphs are provided in Appendix D.

4.3 FORMER MGP STRUCTURES

The test pit and soil boring activities conducted during the Site Characterization confirmed that below-grade MGP structure remnants are present at the Site. Former MGP structures identified in the Manufactured Gas Plant History Report (*GEI*, 2003) are depicted on Figure 2. A summary of below-grade structures encountered during the Site Characterization and RI is provided below.

250,000 Cubic Foot Holder

A total of three test pits (TP-01, TP-02 and TP-03) and three soil borings (SB-01, SB-04 and SB-08) were excavated in the footprint of the former 250,000 cubic foot holder. Based on field observations in these test pits and borings the bottom of the holder was encountered in all the test pits at depths ranging from 8.5 feet below grade (TP-01) to 9.5 feet below grade (TP-03). Similarly, the apparent holder bottom was encountered at depths ranging from approximately 9 feet below grade (SB-04) to 10 feet below grade (SB-01). It should be noted that wood debris was encountered at 15 to 18 feet below grade in SB-01, suggesting a deeper holder bottom.

Remnants of the walls for the former 250,000 c ubic foot holder were encountered at depths ranging from 3.5 to 8.5 feet below grade (TP-02) to 0.5 to 9.5 feet below grade (TP-03). Wood debris was encountered

Groundwater and or NAPL were not detected in any of the test pits or soil boring completed within the former 250,000 holder. Coal tar NAPL was detected approximately 20 feet below grade (i.e., approximately 10 feet below the holder bottom) in soil boring SB-04.

Based on the cross-sections (Figures 7A to 7C), it is estimated that the bottom of the 250,000 cubic foot holder was located at depth of 18 feet below the lower portion of the site.

Retort House

A total of one test pit (TP-02) and one soil boring (SB-12) were excavated in the footprint of the former retort house.

Based on field observations in this test pit and boring remnants of the former retort house were encountered at depths of 4.5 to 6.5 feet bgs. It is likely that the remnants encountered represent the bottom of the structure,

Groundwater was detected in soil boring SB-12 at a depth of 7 feet bgs in soil boring SB-12. Coal tar NAPL was detected at approximately 11.5 feet bgs to 14 feet bgs, and 36 feet bgs to 38 feet bgs in soil boring SB-12.

Based on the cross-sections (Figures 7A to 7C), it is estimated that the bottom of the retorts were located at depth of 14 feet below the lower portion of the Site.

Former Suspected MGP Facility Piping

During Site clearing activities, two 12-inch vertical cast iron pipe openings were uncovered north of test pit TP-4A, adjacent to the concrete block retaining wall, south of the 50,000 cubic foot holder. These pipe openings were observed to be open-ended couplings to the two 12-inch cast iron pipes and are likely to be associated with the onsite former MGP structures, likely the 50,000 cubic foot holder. It is believed these pipes formerly serviced the 50,000 cubic foot holder. NAPL was observed floating on the water. A sample of this NAPL (Exposed Pipe 1) was collected and submitted to META Environmental located in Watertown, MA for fingerprint analyses for MGP related residuals.

A total of two test pits (TP-4A and TP-4B) and one soil boring (SB-06) were excavated near these pipes. During excavation of TP-4A, the two pipes were exposed from 2 to 3 feet bgs. These pipes were observed extending in a southwestern trajectory under a subsurface square brick structure encountered from approximately 0.5 to 2 feet bgs. The pipes were not

encountered in TP-4B. Neither groundwater nor NAPL were encountered within the excavation of TP-4A or TP-4B. Due to the limited depth of these test pits, SB-06 was advanced in this area to identify any impacts related to the two pipes. NAPL was observed from 7 to 9 feet bgs.

Additionally, per NYSDEC's request, test pit TP-06 was excavated west of the railroad tracks along Federal Street in efforts to locate remnants of the "4" NAPHTHA PIPE", asshown on the 1886 S anborn Fire Insurance Map. No evidence of the former 4-inch naphtha pipe was detected during the test pit excavation.

22,800 Cubic Foot Holder

A total of one test pit (TP-05) and one soil boring (SB-05) were excavated in the footprint of the 22,800 cubic foot holder. Based on field observations in the test pit, the holder wall was encountered from 1.5 to 9 feet bgs.

Groundwater was encountered at 9 feet bgs within soil boring (SB-05). NAPL was detected at a depth of 22 feet bgs within soil boring SB-05.

Based on the cross-sections (Figures 7A to 7C), it is estimated that the bottom of the 22,800 cubic foot holder was located at a depth of 20 feet below the lower portion of the site.

50,000 Cubic Foot Holder

During the City of Yonkers' removal of its salt pile from the area of the Site adjacent to the concrete block retaining wall in the northern boundary of the lower (southern) portion of the Site, the southern edge of the former 50,000 cu ft holder wall was uncovered. Approximately 1 foot of this holder wall was exposed and was observed to be extending underneath the concrete block retaining wall. Further excavation could not be attempted in this location due to the presence of the retaining wall.

During excavation of soil boring SB-10, concrete and wood, which are interpreted to be remnants of the bottom of the former holder, were encountered in the 20 to 22 feet bgs interval. NAPL was observed in this boring at approximate depths of 27 feet bgs to 42.5 feet bgs.

Based on cross-sections (Figures 7A to 7C), it is estimated that the bottom of the 50,000 cubic foot holder was located 18 feet below grade of lower portion of the Site.

4.4 ANALYTICAL RESULTS AND FIELD OBSERVATIONS

4.4.1 SUBSURFACE SOIL

A total of 22 subsurface soil samples were collected from the test pit and soil borings as part of the RI. In addition, a total of six QA/QC samples that included four matrix spike/matrix spike duplicate and two field duplicate samples were also collected. Soil samples were submitted to Chemtech and analyzed for TCL VOCs, TCL SVOCs, TAL metals, and cyanide as described in Section 3. The analytical results of the soil samples collected during the SCS and RI are summarized in Table 3 and presented on Figures 4, 5, and 6. The soil sample analytical data in the tables and figures are compared Soil Cleanup Objectives (USCOs) per 6 NYCRR Part 375 (NYSDEC, 2006) for unrestricted site use. Use of unrestricted SCOs is considered to be conservative since the Site is currently and will be used for commercial/industrial purposes in the

forseeable future. The subsurface soil analytical data and field observations of NAPL and total VOCs measured in the field using PID are summarized below.

Headspace VOCs (PID Measurements)

PID readings for soil samples collected during soil boring/monitoring well installations ranged from not detected to 2,254 ppm. The highest total-VOCs concentration measured using the PID of 2,254 ppm was detected in soil sample soil SB-10 (30 - 32 ft bgs). This elevated PID reading is attributable to the presence of residual NAPL observed between 27 and 42.5 feet bgs in this boring. In general, high PID readings were detected in soil associated with residual NAPL and or staining.

NAPL Observations

During the RI, NAPL was observed in soils collected from:

- SB-10 (approximately 27 to 42.5 feet bgs). This boring was located within the footprint of the 50,000 cubic foot holder. NAPL was observed above and below the concrete and brick debris assumed to be related to the holder wall at this location. As shown on Figures 7B and 7C), the NAPL is occurs above the assumed holder bottom;
- SB-12 (approximately 11.5 14 and 36 38 feet bgs). This boring was located within the footprint of the former retort house. NAPL was observed shallow, within a sand layer at the approximate depth of the bottom of the retort house (Figure 7A). NAPL was also observed deeper within this boring, within weathered bedrock;
- SB-13 (approximately 47 to 49 feet bgs). This boring was located down gradient of the 50,000 cubic foot holder. NAPL was observed within weathered bedrock;
- SB-16 (approximately 30 to 36 feet bgs). This boring was located down gradient of the 50,000 c ubic foot holder. NAPL was observed within layers of bricks, presumably related to the former holder; and,
- MW-7 (approximately 35 to 40 feet bgs). This boring was located down gradient of the 50,000 c ubic foot holder. NAPL was observed within a clay layer, corresponding to the approximate depth of the former holder bottom (Figure 7C).

VOCs

A total of 22 VOCs were detected at least once in the soil samples collected during the RI. Of these, seven VOC, namely acetone, benzene, ethyl benzene, methylene chloride, toluene, o-xylene, and m/p-xylene) were detected at concentrations exceeding their respective USCOs. Acetone and methylene chloride are common laboratory artifacts and are not considered as COI.

Benzene was detected above its SCO in four soil samples [MW-7 (36-38), SB-10 (30-32), SB-13 (47-49), SB-16 (30-33)].

Ethylbenzene, toluene, o-xylene, and m/p-xylene were detected at concentrations exceeding their respective SCOs in four soil samples [MW-7 (36-38), SB-10 (30-32), SB-12 (12-14), SB-16

(30-33)]. Total VOCs concentrations in all soil samples ranged from non-detect to 1,986 milligrams/kilogram (mg/kg), which was detected in subsurface soil sample SB-10 (30 -32).

Exceedances of VOCs and or total VOCs were only detected in soil samples from borings where corresponded to the presence of NAPL was observed, namely SB-10, SB-12, SB-13, SB-16 and or SB-MW-7 within the respective soil borings. The vertical extent of VOCs impacts was delineated at each soil boring and test pit sample location (i.e. no USCOs were exceeded in the deepest sample collected), with the exception of locations SB-13 and SB-16.

In soil boring SB-13, benzene was detected above its USCO in the 47-49 ft bgs sample. However, the benzene concentration detected at SB-13 (0.085 mg/kg) was considerably lower than the Commercial SCO of 44 mg/kg. This was the interval directly above the weathered bedrock and contained NAPL.

In soil boring SB-16, acetone, a common laboratory contaminant was detected above its USCO in the 45-47 ft bgs sample. However, the acetone concentration detected at SB-16 (0.071 mg/kg) was considerably lower than the Commercial SCO of 500 mg/kg.

No other VOCs were detected above their respective USCOs in this sample.

SVOCs

A total of 23 SVOCs were detected in soil samples collected during the RI. Of these, fifteen Polycyclic Aromatic Hydrocarbons (PAHs) (acenaphthene, acenaphthylene, anthracene, benzo(a)anthracene, benzo(a)pyrene, benzo(b)fluoranthene, benzo(k)fluoranthene, chrysene, dibenz(a,h)anthracene, fluoranthene, flourene, indeno(1,2,3-cd)pyrene, naphthalene, phenanthrene, and pyrene) were detected at concentrations exceeding their respective USCOs in soil samples MW-7 (36-38), SB-10 (30-32), SB-12 (12-14), and SB-16 (30-33).

Total SVOC concentrations ranged from non-detect to 13,716 mg/kg, which was detected at SB-10 (30-32) in a sample collected at a depth of 30 to 32 feet bgs.

Exceedances of SVOCs and or total SVOCs were only detected in soil samples from borings where NAPL was observed, namely SB-10, SB-12, SB-13, SB-16 and or SB-MW-7. The vertical extent of SVOC impacts were delineated at each soil boring, monitoring well boring, and test pit sample location.

Inorganics

A total of 23 inorganic constituents were detected in soil samples collected during the RI. Of these, seven inorganics (arsenic, chromium, copper, lead, mercury, nickel, and zinc) were detected at concentrations exceeding their respective USCOs in soil samples TP-6 Floor (5), MW-8 (9.5-10), MW-9 (9-9.5), SB-9 (5-8), SB-9 (28-30), SB-10 (30-32), SB-11 (5-5.5).

The vertical extent of impacts from inorganics was delineated at each soil boring and monitoring well boring, with the exception of location SB-9 where nickel was detected above the USCO in the deepest soil sample (collected from depth of 28 to 30 feet bgs). Additionally, inorganic constituents were detected at concentrations above USCOs in the bottom of test pit TP-06.

It should be noted that the majority of inorganic present above their respective USCO's are within soils encountered at MW-8, MW-9, SB-11, and TP-06, which are all located down gradient and off-site from the historical extent of the former MGP site (Figure 6). Sanborn Fire

Insurance Maps indicate that this offsite area was formerly in an area formerly identified as "open water" on historic maps of this area of Yonkers. Accordingly, this entire off-site area consists of fill material, which typically contains inorganics above USCOs. Metals are commonly detected in soil used as fill in urban areas.

4.4.2 GROUNDWATER

Two rounds of groundwater sampling were conducted as part of the RI during November 2012 and April 2013.

During the November 2012 sampling event, six groundwater monitoring wells (MW-1, MW-3, MW-4, MW-7, MW-8, MW-9) were sampled. MW-2 and MW-5 were not sampled during this round because they had been inadvertently destroyed by current site operations. These wells were replaced in December 2012.

During the April 2013 sampling event eight monitoring wells (MW-1, MW-2, MW-3, MW-4, MW-5, MW-7, MW-8, MW-9) were sampled. Groundwater samples were analyzed for TCL VOCs, TCL SVOCs, TAL Metals, and total cyanide. Due to the high turbidity observed at monitoring wells MW-1 and MW-4 during the April 2013 sampling event, unfiltered and laboratory filtered samples for metals analysis were collected from these wells.

Laboratory analytical results for constituents detected in the groundwater samples are summarized in Table 4. A nalytical results were compared with Ambient Water Quality Standards and Guidance Values (AWQSGV) and guidance values contained in NYSDEC Technical and Operational Guidance Series (TOGS) 1.1.1 (NYSDEC, 1998) these standards and guidance values are protective of groundwater quality assuming that groundwater is used as a source of drinking water. That assumption is not applicable to the Site because groundwater is not used as a source of drinking water. Thus, the use of Class GA standards and guidance values for comparison to Site groundwater is conservative. The analytical results of the groundwater samples collected from each well are presented on Figures 9, 10, and 11. Field measurements and observations as well as analytical results from the groundwater investigation are summarized below.

November 2012 Sampling Round

Groundwater analytical results are summarized in Table 4 and on Figures 9, 10, and 11.

Field Measurements

With the exception of monitoring well MW-4, each monitoring well was sampled after purging and water quality parameters stabilized as described in Section 3.7. Monitoring well MW-4 exhibited a very slow recharge rate and ran dry several times prior to stabilization of water quality parameters. Accordingly, due to the small volume of water produced in this well a dissolved metal sample was not collected during this round of sampling.

During groundwater sampling activities, each well was monitored for the presence of NAPL. No NAPL or sheen was detected in any of the wells, with the exception of monitoring well MW-3 where NAPL globules and sheen were observed and monitoring well MW-7 where sheen was observed. Visual descriptions and observations made during the groundwater sampling activities are presented in the groundwater sampling records provided in Appendix C.

VOCs

A total of 12 VOCs were detected at least once in the groundwater samples collected during the November 2012 sampling event.

- Of these, seven VOCs (isopropylbenzene, styrene, benzene, ethylbenzene, toluene, o-xylene, and m/p-xylene) were detected at concentrations exceeding their respective AWQSGVs in groundwater samples collected from MW-3 and MW-7.
- Of the VOCs, the highest concentrations were detected for benzene, which was detected in groundwater monitoring wells MW-3 (NAPL observed within well) and MW-7 (screened through the interval where NAPL was observed in soil from 35 to 40 feet bgs). No VOCs were detected above their AWQSGVs in any other samples collected during the initial round of groundwater sampling.

SVOCs

Nine PAHs and five other SVOCs were detected at least once in the groundwater samples collected during the November 2012 sampling event. Of these, three SVOCs (biphenyl (diphenyl), bis(2-ethylhexyl)phthalate, and naphthalene) were detected at concentrations exceeding their respective AWQSGVs.

- bis(2-ethylhexyl)phthalate is a common laboratory artifact and is not considered to be present. Biphenyl (diphenyl) and naphthalene were detected above their respective AWQSGV in groundwater samples collected from MW-3 and MW-7.
- Naphthalene was the most commonly detected SVOC, was detected above the Class GA AWQSGV in both MW-3 (NAPL observed within well) and MW-7 (screened through the interval where NAPL observed in soil, 35 to 40 feet bgs).

No other SVOCs were detected above their AWQSGVs in any of the monitoring wells during the first round of sampling.

Inorganics

A total of 24 inorganics were detected at least once in groundwater samples collected during the November 2012 sampling event. Of these, eighteen inorganics (antimony, arsenic, barium, beryllium, cadmium, chromium, copper, iron, lead, magnesium, manganese, mercury, nickel, selenium, sodium, thallium, zinc, and cyanide) were detected at concentrations exceeding their respective AWQSGVs.

- Antimony was detected above its AWQSGV in groundwater collected from four monitoring wells (MW-1, MW-3, MW-7, and MW-8).
- Iron was detected above its AWQSGV in groundwater collected from five monitoring wells (MW-1, MW-3, MW-4, MW-8, and MW-9).
- Lead was detected above its AWQSGV in groundwater collected from two monitoring wells (MW-4 and MW-8(dup)).

- Magnesium was detected above its AWQSGV in groundwater collected from three monitoring wells (MW-4, MW-7, and MW-9).
- Manganese was detected above its AWQSGV in groundwater collected from six monitoring wells (MW-1, MW-3, MW-4, MW-7, MW-8, and MW-9).
- Sodium was detected above its AWQSGV in groundwater collected from five monitoring wells (MW-1, MW-3, MW-4, MW-7, and MW-8).
- Monitoring well MW-4 exhibited an additional twelve inorganic constituents (arsenic, barium, beryllium, cadmium, chromium, copper, mercury, nickel, selenium, thallium, zinc, and cyanide) which were detected above their respective AWQSGVs.
- It is noted that of these; iron, manganese, magnesium and sodium are naturally-occurring inorganics commonly detected in groundwater. A dditionally, many other metals are commonly present in rock salt including lead and zinc. Iron cyanide is also often added to rock salt used for roadway de-icing as an anti-caking agent. The presence of many these inorganics in groundwater appear to be related to the storage of road salts at the lower portion of the site or in the case of offsite monitoring wells MW-8 and MW-9, the presence of historic fill containing these analytes.

April 2013 Sampling Round

Groundwater analytical results are summarized in Table 4 and on Figures 9, 10, and 11.

Field Measurements

Each monitoring well was sampled upon reaching parameter stability and turbidity levels below 50 NTU with the exception of MW-1 and MW-4; both had very slow recharge and ran dry several times prior to stabilization of water quality parameters. A sample for dissolved metals was collected from both wells.

During groundwater sampling activities, each monitoring well was monitored for the presence of NAPL. No NAPL or sheens were noted in any of the wells with the exception of MW-3 and MW-7 where NAPL globules and sheens were observed. Visual descriptions and observations made during the groundwater sampling activities are presented on the groundwater sampling records provided in Appendix C.

VOCs

A total of 14 VOCs were detected at least once in the groundwater samples collected during the April 2013 sampling event.

- Of these, eight (8) VOCs (acetone, isopropylbenzene, styrene, benzene, ethylbenzene, toluene, o-xylene, and m/p-xylene) were detected at concentrations exceeding their respective AWQSGVs.
- Isopropylbenzene was detected above its AWQSGV in groundwater collected from monitoring well MW-7.

- Styrene was detected above its AWQSGV in groundwater collected from monitoring well MW-3.
- Benzene, ethylbenzene, toluene, o-xylene, and m/p-xylene were detected above their respective AWQSGVs in three (3) monitoring wells (MW-3, MW-5, and MW-7).

No VOCs were detected above their AWQSGVs in any other sampled monitoring wells.

SVOCs

Thirteen PAHs and six other SVOCs were detected at least once in the groundwater samples collected during the April 2013 s ampling event. O f these, nine (9) were detected at concentrations exceeding their AWQSGVs (biphenyl (diphenyl), acenaphthene, benzo(a)anthracene, benzo(a)pyrene,benzo(b)fluoranthene, chrysene, fluorene, naphthalene, and phenanthrene).

- Biphenyl (diphenyl), acenaphthene, benzo (a) anthracene, benzo (b) fluoranthene, chrysene, and naphthalene were detected above their respective AWQSGVs in three (3) monitoring wells (MW-3, MW-5, and MW-7).
- Fluorene was detected above its AWQSGV in the duplicate groundwater sample collected from monitoring well MW-3.
- Phenanthrene was detected above its AWQSGV in groundwater collected from monitoring wells MW-3 and MW-7.

No other SVOCs were detected above their AWQSGVs in any of the monitoring wells.

Inorganics

A total of 23 inorganics were detected at least once in groundwater samples collected during the April 2013 sampling event. Of these, 16 inorganics (antimony, arsenic, barium, beryllium, chromium, copper, iron, lead, magnesium, manganese, mercury, nickel, selenium, sodium, thallium, and cyanide) were detected at concentrations exceeding their respective AWQSGVs.

- Antimony was detected above its AWQSGV in groundwater collected from two monitoring wells (MW-2 and MW-5).
- Arsenic, chromium, copper, mercury, nickel, selenium, and thallium were detected above their respective AWQSGV s in groundwater collected from monitoring well MW-4.
- Barium was detected above its AWQSGV in groundwater collected from monitoring well MW-2.
- Iron, manganese and sodium were detected above their respective AWQSGVs in each of the seven monitoring wells sampled (MW-1, MW-2, MW-3, MW-4, MW-5, MW-7, MW-8, and MW-9).

- Lead was detected above its AWQSGV in groundwater collected from three monitoring wells (MW-3, MW-4, and MW-5).
- Magnesium was detected above its AWQSGV in groundwater collected from three monitoring wells (MW-5, MW-7, and MW-9).
- Thallium was detected above its AWQSGV in groundwater collected from monitoring well MW-9.
- Cyanide was detected above its AWQSGV in groundwater collected from monitoring well MW-2.

No other inorganics were detected above their AWQSGVs in any of the monitoring wells.

A total of 15 dissolved inorganics were detected at least once in groundwater samples collected during the RI. Of these, four inorganics (antimony, manganese, selenium, and sodium) were detected at concentrations exceeding their respective AWQSGVs. Manganese and sodium were detected above their respective AWQSGVs in groundwater collected from monitoring well MW-1 (up gradient well). Antimony, manganese, selenium, and sodium were detected above their respective AWQSGVs in groundwater collected from monitoring well MW-4. Dissolved organics sampling was limited to MW-1 and MW-4.

4.4.3 SOIL GAS SAMPLE RESULTS

The NYSDOH Guidance does not provide screening values for VOCs in soil vapor for comparison. A total of 15 VOCs were detected at least once in the soil gas samples collected during the RI. The detected compounds are characterized into three general categories: chlorinated VOCs (CVOCs), non-MGP related compounds, and possibly MGP related compounds.

CVOCs

1,1,1-Trichloroethane (TCA) and tetrachloroethene (PCE) were the only two CVOCs detected in the soil gas samples collected from the site. PCE was detected in all three samples, with the highest concentration present in SV-3. CVOCs are not soil gas contaminants associated with former MGP sites. TCA is a common degradation byproduct of PCE.

Non-MGP Related Compounds

Nine VOCs that are not considered MGP related or CVOCs were detected at low concentrations in the soil gas samples collected from the site, including 1,3-butadiene, 2,2,4-trimethylpentane, 2-butanone (methyl ethyl ketone), acetone, carbon disulfide, cyclohexane, Freon 12 (Dichlorodifluoromethane), heptanes, and hexane. The greatest total concentration of non-MGP related compounds was detected in SV-4. Many of these VOCs can be lab artifacts.

Possibly MGP Related Compounds

Compounds potentially related to MGP residues detected in the soil gas samples include benzene, toluene, ethylbenzene, and xylenes (BTEX). The greatest total concentration of possibly MGP related compounds was detected in SV-4. However, a review of data collected during SC and RI activities from MW-4 and SB-02 indicates the absence of these analytes in soil

or groundwater, beyond an estimated concentration of benzene within soil at MW-04. Additionally, a review of the boring logs from both of these locations indicates the absence of any NAPL, staining, odors, or elevated PID readings, suggesting that the detection of these analytes within soil gas at SV-4 is from a non-MGP related source.

EXTENT OF MGP NAPL IMPACTS

As noted in earlier sections of this document, the surface elevation of the lower and upper portion differ greatly, ranging from a surface elevation of 16.9 feet MSL at SB-04 to a surface elevation of 39.7 feet MSL at SB-13. The approximately 23 feet variance in grade is due to the City of Yonkers' placement there of various debris, across the northern portion of the Site and is the reason for its designation as the 'upper' portion of the Site.

In an effort to effectively show the distribution of coal tar residue in the subsurface, Figures 12 to 15 the occurrence of NAPL relative to remnant and former MGP structures and depth in the subsurface relative to the ground surface of the lower portion for the site are presented as depths relative to ground surface of the lower portion of the Site. For example, Figure 12 depicts the location of where NAPL was observed 6 to 11 feet below the assumed average ground surface of the lower portion of the Site (16 feet msl).

5.1 MGP NAPL - 6 TO 11 FEET

Figure 12 presents the location of relatively shallow NAPL detected on the Site at 6 to top of water table (approximately 11 feet below grade). This plan view NAPL area is approximately 50 feet wide and 100 feet long or 5,000 square feet. The northern section of this NAPL area, from the concrete block retaining wall towards SB-10 and SB-16 is likely related to the former 50,000 cubic foot holder. Cross-sections appear to indicate that the southern portion of this area is occurs down gradient of the former 22,800 cubic foot holder and the former 50,000 foot holder. As discussed previously, the VOC and SVOC impacts within soil are concurrent with and related to the presence of NAPL.

It should be noted that the area immediately to the north of the concrete block retaining wall was inaccessible due to a steep unstable slope. While the two NAPL areas are presumed to be connected as shown, it is possible that the areas are separate. This portion of NAPL impacts, while shown on this figure as present 6 to 11 feet below the lower parcel, some of these areas are covered with an additional 23 feet of fill. It is anticipated that this area will be investigated further during a pre-design investigation (PDI).

5.2 MGP NAPL – 11 TO 21 FEET

Figure 13 presents the occurrence of NAPL detected between 11 to 21 feet below grade. This area is approximately equivalent to a rectangle 75 feet wide and 100 feet long, or 7,500 square feet. Cross-sections on Figure 7A and Figure 7B indicate that the western section of this NAPL area at SB-07 and SB-12 are from the former 250,000 cubic foot holder or from the down gradient edge of the NAPL from the two smaller former gas holders. NAPL observed from the concrete block retaining wall towards SB-10 and SB-16 is related to the former 50,000 cubic foot holder. NAPL at SB-05 and SB-06, the southern portion of this area, is present down gradient of the former 22,800 cubic foot holder and the former 50,000 foot holder. As discussed previously, the VOC and SVOC impacts within soil are concurrent with and related to the presence of NAPL.

It should be noted that the area between SB-12 and MW-7, as well as the area immediately to the west of MW-7, were inaccessible due to a steep unstable slope. It is anticipated that this area will be investigated further during a pre-design investigation (PDI).

5.3 MGP NAPL – 21 TO 25 FEET

Figure 14 depicts the location of NAPL detected on the Site at 21 to 25 feet below the surface of the lower parcel. The NAPL detected at these depths are isolated at SB-04 (250,000 cubic foot gas holder), SB-05 (22,800 cubic foot gas holder), and SB-13 (50,000 cubic foot gas holder). The NAPL at SB-05 and SB-13 at this depth, appear to be the lower edges of the NAPL noted in Section 5.4. As discussed previously, the VOC and SVOC impacts within soil are concurrent with and related to the presence of NAPL.

5.4 MGP NAPL – DEEPER THAN 35 38 FEET

Figure 15 depicts the location of NAPL detected on the Site deeper than 35 feet below grade. The NAPL detected at this depth is isolated to the vicinity of soil boring SB-12, at a depth above bedrock. It is believed that this NAPL migrated from upgradient locations at the two smaller holders. NAPL was not detected at SB-11, the down gradient location of boring SB-12. As discussed previously, the VOC and SVOC impacts within soil are concurrent with and related to the presence of NAPL.

EXPOSURE ASSESSMENT

Information collected during the initial Site Characterization and the RI at the former Ludlow Street Works Site has been used to qualitatively assess potential exposure pathways for the various detected compounds in Site soils, groundwater, and soil vapor. The Site is located within an industrial district of a multi-use (i.e., residential and commercial) neighborhood and is currently used as a City of Yonkers Department of Public Works storage yard (lower portion of the Site) and fill deposition area (upper portion of the Site)

6.1 SOILS – VOCS, SVOCS AND MGP NAPL

As noted in earlier sections of this document, NAPL was detected at discrete locations in the subsurface ranging in depths from 6 to 38 feet below the surface of the lower portion of the Site. One VOC was detected above its USCO as shallow as 3 feet below ground surface (toluene at TP-4). However, all other VOC and SVOC impacts were concurrent with the observations of NAPL at onsite areas, with the highest VOC and SVOCs detected associated with NAPL in the vicinity of the foundation of the former 50,000 cubic foot gas holder. As discussed in Section 5, the majority of NAPL is present at depths deeper than 11 feet below the lower portion of the Site, and is covered with up to 40 feet of fill materials.

There is limited potential for shallow NAPL impacted soil to be encountered during intrusive maintenance activities adjacent to the concrete block wall. It is unlikely that these materials would be encountered during typical Site operations (i.e., use as a DPW storage yard). Most Site soils are covered with gravel and broken bituminous pavement and, for most of the year; the southern (or lower) portion of the Site is covered by a pile of road salt, limiting the potential exposure to humans or the environment.

VOCs, SVOCs and MGP NAPL are not present offsite, west of the railroad tracks.

6.2 SOILS – INORGANIC IMPACTS

Seven inorganics (arsenic, chromium, copper, lead, mercury, nickel, and zinc) were detected at concentrations exceeding their respective USCOs. As noted in earlier portions of this document, the majority of inorganics present above USCO are within offsite soils encountered at MW-8, MW-9, SB-11, and TP-06, locations down gradient and off-site from the historical extent of the former MGP site. Sanborn Fire Insurance Maps indicate that this offsite area was formerly in an area formerly identified as open water. Accordingly, this entire off-site area consists of non-native urban fill, which typically contains metals and other inorganic potential COI at concentrations above their SCOs.

Given the current depth below ground surface of remnant MGP structures, and the presence of the fill emplaced on the upper portion of the Site, it is assumed that these areas have been covered by historic fill materials, containing metals and other inorganics. As noted in Section 6.1, the lower portion of the site is covered by pavement, soil, and a salt pile, while the upper portion of the Site is covered approximately 15 feet of construction debris (e.g., wood, concrete, boulders, asphalt, gravel, etc.) and some vegetation. These cover materials limit the potential exposure to humans or the environment from the historic fill materials.

6.3 GROUNDWATER - VOCS, SVOCS AND MGP NAPL

Analytical results indicated the presence of VOCs, PAHs, and inorganic concentrations in groundwater at the Site above the Class GA AWQSGV and guidance values. Two monitoring wells (MW-3 and MW-7) exceeded AWQSGV MGP-related VOCs and SVOCs during the first groundwater sampling round, and again during the second groundwater sampling round, along with replacement well MW-5.

Groundwater at the Site is currently not used for a potable water source and there are no plans for future use of potable or commercial/industrial groundwater at the Site. Groundwater flow direction is predominantly to the west - northwest towards the Hudson River. The depths to groundwater measured in on and off-site wells in January 2013 ranged from approximately 4 to 39 feet bgs which corresponds to elevations of approximately 2 to 6 feet MSL. Therefore, the potential for exposure to groundwater during intrusive subsurface activities is not anticipated.

6.3 GROUNDWATER – FILL RELATED CONSTITUENTS

Several inorganic analytes were detected at the Site (including sodium, magnesium, and manganese) at concentrations exceeding Class GA Groundwater Standards. These analytes are not associated with MGP related operations. Their presence at high concentrations in both Site soils and in Site groundwater is believed to be due to use of the Site for road salt storage, or due to the presence of urban fill materials. These analytes are also present in offsite groundwater, down gradient of the Site. The presence of these analytes in the off-site downgradient wells is consistent with their proximity to the Hudson River, which is a saline water body.

6.4 SOIL VAPOR

Soil vapor results indicate the presence of 15 VOCs in the soil gas samples collected during the RI. In an effort to assess the potential impact of these concentrations, these concentrations were compared to target concentrations in the document "USEPA OSWER Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance)" Table 2b, EPA530-D-02-004, November 2002. As noted on Table 5, none of the possibly related MGP VOCs exceeded the Target Soil Gas Concentrations. Based on this comparison, there appears to be no exposure risk to offsite buildings, related to soil vapor concentrations from MGP related contaminants.

CONCLUSIONS AND RECOMMENDATIONS

7.1 CONCLUSIONS

The results of the RI and SCS show that impacted soil and groundwater are associated with discrete on-site source areas in the southern portion of the Site where the former works operations took place. Due to the depth and occurrence of the impacted media and configuration of the Site, there are no current exposure risks. Impacts detected on the northern portion are isolated to the southern-most areas in the vicinity of the former 50,000 cu ft gas holder and retort house. The following provides a summary of the key findings.

- Remnants of the former MGP structures (lime house, purifying house, meter house, retort house 250,000 cubic foot holder, 22,00 cubic foot holder, 50,000 gas holder are present on the property.
- No evidence of the former tar pipe identified on the 1886 Sanborn Fire Insurance maps was observed in the field.
- An active sanitary force main is present offsite (21" vitrified clay east of Site); and abandoned sewer main (14-inch cast iron pipe) are present in the subsurface at the Site and offsite areas. The force main is operated by the Westchester County Department of Environmental Facilities. This active conveyance is present at a depth of approximately 12 feet bgs relative to the grade at the southern or lower portion of the Site and runs north-south along the eastern border of the Site. The 14-inch cast iron pipe is abandoned in place, generally bisecting the Site east-west through the northern portion of the Site, from the sanitary force main located along the eastern boundary of the Site.
- A storm sewer main, constructed of reinforced concrete pipe, according to the City of Yonkers, runs to the west-northwest from the southeast corner of the southern portion of the Site. This structure appears to extend beneath the MNRR right-of-way.
- Surface soil across the southern portion of the Site, where the predominance of MGP
 activities occurred, is covered by road salt, fill material, isolated remnant asphalt areas
 and abandoned concrete structures. Exposure to surface soils in the southern portion of
 the site is limited due to the extensive fill, salt piles and controlled access.
- Surface soil in the northern portions of the site that was present at grade when the MGP was operational, is currently covered by 15 to 20 feet of debris and fill that was placed there by the City of Yonkers subsequent to operations of the former MGP. Accordingly, there is no potential to encounter the former surface soils there. Furthermore, based on the historical records, this portion of the Site was not utilized during MGP operations and so it is not expected that former surface soils there contain impacts from the former MGP.

- Subsurface soil is comprised of three unconsolidated units, which are fill, sand, silt and clay, with isolated deposits of sand and cobbles, and gravel and sand, as shown on the cross sections on Figure 7A to Figure 7C. Fill, consisting of non-native soils, bricks, concrete and other debris, is encountered to depths up to 14 feet below the lower portion of the Site, below which sand is encountered to bedrock. In some areas of the site, the silt and clay layer is encountered from 14 to 24 feet below the lower portion of the site. NAPL and groundwater are predominantly found in the sand layer. Ground water is encountered primarily within the sand layer, at an approximate depth of 14 feet below the lower portion of the site.
- Bedrock occurs at depths ranging from approximately 20 to 36 feet below grade of the lower portion of the site.
- Subsurface soil at the southern portion of the Site is impacted by residual coal tar and NAPL. The soil also contains elevated concentrations of VOCs and SVOCs associated with the 22,800 cu foot holder, the 250,000 cu foot holder, and the retort house (Figure 12 to Figure 15). The impacts in the areas surrounding these structures occur primarily in the soils within the respective structures and deeper soils outside of and in the vicinity of the structures above and below the water table, with the shallowest occurrences observed at depths below 6 to 11 feet bgs below the lower portion of the Site.
- Subsurface soil in the southern-most area of the northern portion of the Site are impacted by residual coal tar and NAPL associated with the 50,000 cu ft holder and retort house (Figure 12 to Figure 15). The impacts occur primarily inside the remnants of these structures and in the soils below the water table to a depth of approximately 25 feet bgs below the lower portion of the Site.
- Impacts were observed in the deep subsurface soil at 36 feet bgs below the lower portion of the Site at SB-12, within the footprint of former retort house and at 25 feet bgs below the lower portion of the site, within weathered bedrock at SB-13, outside of the 50,000 cu foot holder
- Based on the historical records, the northern portion of the Site (north of the footprint of the 50,000 cubic foot holder), was not utilized during MGP operations and accordingly it is not likely that subsurface soils in this area contain impacts related to the former MGP.
- Groundwater occurs at depths that ranged from 9 to 14 feet bgs of the lower portion of the Site and flows from the southeast to northwest across the Site (Figure 7). Where groundwater flows through soil containing MGP source material (e.g., coal tar NAPL or residual coal tar) it leaches the more soluble components of the coal tar, which include VOCs and SVOCs. As a result, groundwater in the vicinity of the source areas (as identified above) contained elevated concentrations of BTEX (VOCs) and some PAHs (SVOCs). This distribution is shown on Figure 9 and Figure 10. As previously noted, it is likely that impacted groundwater has migrated down gradient from the site to areas beneath the railroad right-of-way. D ue to the mild influence of tidal fluctuation on groundwater flow in areas closer to the Hudson River and the numerous natural

attenuation processes that occur in the subsurface (i.e., dilutions, dispersion, adsorption, biodegradation, etc.) it is not likely that impacted groundwater migrates a substantial distance from the on-site source areas.

- VOCs were detected at low levels in soil gas samples collected beneath buildings that currently occupy parcels adjacent to the site. There was no apparent evidence that VOCs detected in soil gas were migrating into the structures at these locations. Therefore, additional monitoring of soil gas is not warranted.
- Elevated concentrations of sodium, manganese, and magnesium detected in Site soil and groundwater are believed to be associated with use of the Site as a road salt storage facility, or due to the presence of urban fill materials.

7.2 **RECOMMENDATIONS**

Based on the RI activities, no additional investigation is recommended to delineate the impacts identified within the Site.

It is recommended that remedial alternatives be evaluated in order to address the subsurface impacts identified during the remedial investigation. P rior to identifying overall remedial approach and specific alternatives, Con Edison will consult with NYSDEC to review and consider the impacts and Site-specific conditions.

SECTION 8

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Table 1 RI Sample Summary

Former Ludlow Street Works Site

Consolidated Edison Company of New York Remedial Investigation - October 2012 through April 2013

Location	Sample ID	Depth (bgs)	TCL VOCs	TCL SVOCs	TAL Metals	Cyanide	TO-15 Modified	Dissolved Metals
		OIL SAMPLES			1		1	
MW-7	MW-7 (36-38)	36-38'	X	X	X	X		
	MW-7 (48-50)	48-50'	X	X	X	X		
) WILL O	MW-17 (36-38)*	36-38'	X	X	X	X		
MW-8	MW-8 (9.5-10)	9.5-10'	X	X	X	X		
) (TY) ()	MW-8 24.5-25)	24.5-25'	X	X	X	X		
MW-9	MW-9 (35-35.5)	35-35.5'	X	X	X	X		
ap 0	MW-9 (9-9.5)	9-9.5'	X	X	X	X		
SB-9	SB-9 (5-8)	5-8'	X	X	X	X		
GD 10	SB-9 (28-30)	28-30'	X	X	X	X		
SB-10	SB-10 (30-32)	30-32'	X	X	X	X		
gp 44	SB-10 (45-47)	45-47'	X	X	X	X		
SB-11	SB-11 (5-5.5)	5-5.5'	X	X	X	X		
GD 10	SB-11 (29.5-30)	29.5-30'	X	X	X	X		
SB-12	SB-12 (12-14)	12-14'	X	X	X	X		
GD 10	SB-12 (36-38)	36-38'	X	X	X	X		
SB-13	SB-13 (47-49)	47-49'	X	X	X	X		
SB-14	SB-14 (38-40)	38-40'	X	X	X	X		
	SB-14 (45-47)	45-47'	X	X	X	X		
~~	SB-14A (45-47)*	45-47'	X	X	X	X		
SB-16	SB-16 (30-33)	30-33'	X	X	X	X		
~~	SB-16 (45-47)	45-47'	X	X	X	X		
SB-17	SB-17 (33-35)	33-35'	X	X	X	X		
	SB-17 (45-47)	45-47'	X	X	X	X		
TP-6	TP-6 Floor	6'	X	X	X	X		
) (TY 1	GROUNDWATE		T		_	т —	1	
MW-1	MW-1	NA	X	X	X	X		
MW-3	MW-3	NA	X	X	X	X		
MW-4	MW-4	NA	X	X	X	X		
MW-7	MW-7	NA	X	X	X	X		
MW-8	MW-8	NA	X	X	X	X		
MW-8	MW-18*	NA	X	X	X	X		
MW-9	MW-9	NA NA	X	X	X	X		
MXX 1	GROUNDWA'	_	_	-		1 37		W
MW-1	MW-1	NA	X	X	X	X		X
MW-2	MW-2	NA NA	X	X	X	X		
MW-3	MW-3	NA NA	X	X	X	X		
MW-3	MW-3A*	NA NA	X	X	X	X		37
MW-4	MW-4	NA NA	X	X	X	X		X
MW-5	MW-5	NA NA	X	X	X	X		
MW-7 MW-8	MW-7 MW-8	NA NA	X	X	X	X		
1 IV I VV _ X	IIVIW-X	I NA	ιX	X	X	ιX	1	

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Table 1

RI Sample Summary Former Ludlow Street Works Site

Consolidated Edison Company of New York Remedial Investigation - October 2012 through April 2013

Location	Sample ID	Depth (bgs)	TCL VOCs	TCL SVOCs	TAL Metals	Cyanide	TO-15 Modified	Dissolved Metals
	SOII	GAS SAMPI	LES					
SV-1/SB-9	SV-1	7'					X	
SV-3	SV-3	8'					X	
SV-4	SV-4	8'					X	

X - Indicates sample was analyzed

^{* -} Indicates a duplicate sample.

Table 2 Summary of Groundwater Elevations Former Ludlow Street Works Consolidated Edison Company of New York Remedial Investigation - October 2012 through April 2013

	Top of Casing	11/19	0/2012	1/3/2	2013	4/5/2	2013
Well ID	Elevation (feet	Depth to	Groundwater	Depth to	Groundwater	Depth to	Groundwater
Well ID	AMSL)	Water	Elevation	Water	Elevation	Water	Elevation
	AWISL)	(feet)	(feet AMSL)	(feet)	(feet AMSL)	(feet)	(feet AMSL)
MW-1	20.35	13.95	6.40	14.06	6.29	13.36	6.99
MW-2	16.24	$NA^{(1)}$	NA ⁽¹⁾	10.45	5.79	10.92	5.32
MW-3	19.35	16.33	3.02	16.51	2.84	16.49	2.86
MW-4	19.28	14.09	5.19	14.15	5.13	13.92	5.36
MW-5	15.16	$NA^{(1)}$	NA ⁽¹⁾	11.23	3.93	12.05	3.11
MW-7	42.03	39.08	2.95	39.19	2.84	39.00	3.03
MW-8	5.20	5.06	2.08	5.20	1.94	NA ⁽²⁾	NA ⁽²⁾
MW-9	6.35	4.23	2.12	4.38	1.97	4.15	2.20

Notes:

- (1) Monitoring well destroyed by DPW site activities, replaced in December 2012.
- (2) Access blocked during sampling activities.

AMSL = Above Mean Sea Level

Elevations are based on the North American Vertical Datum of 1988 (NAVD88).

					Dup of MW-7(36-38)					
Con Ed - Luc			Location ID:	MW- 7	MW- 7	MW- 7	MW-8	MW-8	MW-9	MW-9
	il Analytical Data		Sample ID:	MW-7(36-38)	MW-17(36-38)	MW-7(48-50)	MW-8(9.5-10)	MW-8(24.5-25)	MW-9(9-9.5)	MW-9(35-35.5)
Detected Cor	mpound Summary		Lab Sample Id	D4710-03	D4710-09	D4710-04	D4751-03	D4751-04	D4751-06	D4751-05
			Depth:	36 - 38 ft	36 - 38 ft	48 - 50 ft	9.5 - 10 ft	24.5 - 25 ft	9 - 9.5 ft	35 - 35.5 ft
			Source:	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH
			SDG:	D4710	D4710	D4710	D4751	D4751	D4751	D4751
		Unrestricted Use	Matrix:	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
		Soil Cleanup	Sampled:	10/26/2012 9:10	10/26/2012 9:15	10/26/2012 9:20	11/6/2012 14:15	11/6/2012 14:20	11/6/2012 10:00	11/6/2012 10:05
		Objectives	Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND		UNITS:							
	VOLATILES									
67-64-1	ACETONE	0.05	mg/kg	ND	ND	ND	0.051	ND	ND	ND
71-43-2	BENZENE	0.06	mg/kg	7.1 J	5.3	0.0013 J	0.0023 J	ND	ND	ND
75-15-0	CARBON DISULFIDE		mg/kg	ND	ND	0.0013 J	0.0049 J	ND	ND	ND
74-87-3	CHLOROMETHANE		mg/kg	ND	ND	ND	ND	ND	ND	ND
110-82-7	CYCLOHEXANE		mg/kg	0.062 J	0.0082 J	ND	ND	ND	ND	ND
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE		mg/kg	ND	ND	ND	ND	ND	ND	ND
95-50-1	1,2-DICHLOROBENZENE	1.1	mg/kg	ND	ND	ND	ND	ND	ND	ND
541-73-1	1,3-DICHLOROBENZENE	2.4	mg/kg	ND	ND	ND	ND	ND	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND	ND	ND	ND	ND	ND
75-71-8	DICHLORODIFLUOROMETHANE		mg/kg	ND	ND	ND	0.0034 J	ND	ND	ND
75-34-3	1,1-DICHLOROETHANE	0.27	mg/kg	ND	ND	ND	ND	ND	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	30 J	48	0.0017 J	ND	ND	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)		mg/kg	6.2 J	3.6	ND	ND	ND	ND	ND
108-87-2	METHYLCYCLOHEXANE		mg/kg	0.21 J	0.046 J	ND	ND	ND	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	ND	ND	ND	ND	ND	ND	ND
100-42-5	STYRENE		mg/kg	0.89	0.96 J	ND	ND	ND	ND	ND
79-34-5	1,1,2,2-TETRACHLOROETHANE		mg/kg	ND	ND	ND	ND	ND	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	15 J	18	0.0013 J	ND	ND	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE		mg/kg	ND	ND	ND	ND	ND	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE		mg/kg	ND	ND	ND	ND	ND	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	34 J	54	0.002 J	ND	ND	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE	0.26	mg/kg	18 J	28	ND	ND	ND	ND	ND
	Total VOCs	NS	mg/kg	111.462	157.9142	0.0076	0.0616	ND	ND	ND
	SEMIVOLATILES						****			
92-52-4	BIPHENYL (DIPHENYL)		mg/kg	7	11	ND	ND	ND	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE		mg/kg	ND	ND	ND	ND	ND	ND	ND
86-74-8	CARBAZOLE		mg/kg	0.55	0.7	ND	ND	ND	ND	ND
132-64-9	DIBENZOFURAN		mg/kg	2.2	2.8	ND	ND	ND	ND	ND
131-11-3	DIMETHYL PHTHALATE		mg/kg	0.221 J	0.226 J	ND	0.215 J	0.221 J	0.292 J	0.2 J
84-74-2	DI-N-BUTYL PHTHALATE		mg/kg	ND	ND	ND	ND	ND	ND	ND
	PAHs									
83-32-9	ACENAPHTHENE	20	mg/kg	15 J	21 J	ND	ND	ND	ND	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	12	16	ND	ND	ND	ND	ND
120-12-7	ANTHRACENE	100	mg/kg	13	20	ND	ND	ND	ND	ND
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	8.3	13	ND	0.26 J	ND	ND	ND

					Dup of MW-7(36-38)					
Con Ed - Lud	llow Site		Location ID:	MW-7	MW- 7	MW- 7	MW-8	MW-8	MW-9	MW-9
Validated So	il Analytical Data		Sample ID:	MW-7(36-38)	MW-17(36-38)	MW-7(48-50)	MW-8(9.5-10)	MW-8(24.5-25)	MW-9(9-9.5)	MW-9(35-35.5)
Detected Cor	npound Summary		Lab Sample Id	D4710-03	D4710-09	D4710-04	D4751-03	D4751-04	D4751-06	D4751-05
			Depth:	36 - 38 ft	36 - 38 ft	48 - 50 ft	9.5 - 10 ft	24.5 - 25 ft	9 - 9.5 ft	35 - 35.5 ft
			Source:	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH
			SDG:	D4710	D4710	D4710	D4751	D4751	D4751	D4751
		Unrestricted Use	Matrix:	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
		Soil Cleanup	Sampled:	10/26/2012 9:10	10/26/2012 9:15	10/26/2012 9:20	11/6/2012 14:15	11/6/2012 14:20	11/6/2012 10:00	11/6/2012 10:05
		Objectives	Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	0.0,111.11	UNITS:		.,,,					
	VOLATILES									
50-32-8	BENZO(A)PYRENE	1	mg/kg	5.7 J	9 Ј	ND	0.26 J	ND	ND	ND
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	4.4 J	6.6 J	ND	0.3 J	ND	ND	ND
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	2.4 J	2.9 J	ND	0.22 J	ND	ND	ND
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	1.7	2.2	ND	ND	ND	ND	ND
218-01-9	CHRYSENE	1	mg/kg	8	12	ND	0.29 J	ND	ND	ND
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	0.72	0.88	ND	ND	ND	ND	ND
206-44-0	FLUORANTHENE	100	mg/kg	14	23	ND	0.57	ND	ND	ND
86-73-7	FLUORENE	30	mg/kg	22	27	ND	ND	ND	ND	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	2.1	2.7	ND	0.17 J	ND	ND	ND
91-57-6	2-METHYLNAPHTHALENE		mg/kg	110	130	ND	ND	ND	ND	ND
91-20-3	NAPHTHALENE	12	mg/kg	220	240	ND	ND	ND	ND	ND
85-01-8	PHENANTHRENE	100	mg/kg	60	79	ND	0.48	ND	ND	ND
129-00-0	PYRENE	100	mg/kg	26	34 J	ND	0.55	ND	ND	ND
	Total PAHs		mg/kg	525.32	639.28	ND	3.1	ND	ND	ND
	Total SVOCs		mg/kg	535.291	654.006	ND	3.315	0.221	0.292	0.2
	INORGANICS									
7429-90-5	ALUMINUM		mg/kg	7150	6870	2050	4920 J	8550 J	13400 J	2710 J
7440-38-2	ARSENIC	13	mg/kg	2.45	2.47	1.55	9.24 J	2.16 J	ND	0.75 J
7440-39-3	BARIUM	350	mg/kg	57.6	54	36.9	73 J	85.6 J	177 J	60.2 J
7440-41-7	BERYLLIUM	7.2	mg/kg	0.28	0.29	0.08 J	ND	0.33	0.13 J	0.07 J
7440-43-9	CADMIUM	2.5	mg/kg	0.48	0.45	0.03 J	1.32	0.66	2.17	0.19
7440-70-2	CALCIUM		mg/kg	1350	1340	709	3220 J	1870 J	2660 J	13000 J
7440-47-3	CHROMIUM, TOTAL	30	mg/kg	13.6 J	13.7 J	4.91 J	31.2 J	15.8 J	28.1 J	6.01 J
7440-48-4	COBALT		mg/kg	7.95	7.44	2.66	6.63	10.7	27.8	3.85
7440-50-8	COPPER	50	mg/kg	16 J	16.2 J	8.56 J	93.7 Ј	18.8 J	7.84 J	8.4 J
7439-89-6	IRON		mg/kg	22400	21300	5950	33500	24400	41900	6990
7439-92-1	LEAD	63	mg/kg	14.4	14.6	3.07	170	16.1	17.4	4.01
7439-95-4	MAGNESIUM		mg/kg	3170	3100	1310	2310 J	4040 J	7840 J	6590 J
7439-96-5	MANGANESE	1600	mg/kg	180	135	44.7	240 J	427 J	251 J	192 J
7439-97-6	MERCURY	0.18	mg/kg	0.011 J	0.006 J	0.003 J	4.03	ND	0.425	ND
7440-02-0	NICKEL	30	mg/kg	23	22	8.83	19.4	26.5	48.6	10.6
7440-09-7	POTASSIUM		mg/kg	1980	1960	562	1530 J	2600 J	10700 J	860 J
7782-49-2	SELENIUM	3.9	mg/kg	ND	ND	0.52	0.57	0.26 J	0.5	ND
7440-22-4	SILVER	2	mg/kg	0.5	0.47	0.1 J	0.36	0.57	0.23 J	ND
7440-23-5	SODIUM		mg/kg	2090	2190	195	627 J	662 J	402 J	1080 J
7440-28-0	THALLIUM		mg/kg	0.46 J	0.48 J	ND	1.23	0.52 J	ND	ND
7440-62-2	VANADIUM		mg/kg	18	17.7	5.98	15.3 J	21.2 J	38.8 J	9.23 J
7440-66-6	ZINC	109	mg/kg	36.1	39.3	13.7	96.8	47.2	97.6	14.5
57-12-5	CYANIDE	27	mg/kg	0.29 J	0.362	0.171 J	0.097 J	ND	0.087 J	0.075 J

Notes:

- (1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)
- (2) -- indicates no cleanup objective or background level is available.
- (3) ND indicates compound was not detected.
- (4) J indicates an estimated concentration.
- (5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.
- (6) NA indicates compound was not analyzed.

Table 3
Summary of Soil Analytical Data
Former Ludlow Street Works Site
Consolidated Edison Company of New York
Remedial Investigation - October 2012 through April 2013

Con Ed - Luc	low Site		Location ID:	SB- 9	SB- 9	SB-10	SB-10	SB-11	SB-11	SB-12
Validated So	il Analytical Data		Sample ID:	SB-9(5-8)	SB-9(28-30)	SB-10(30-32)	SB-10(45-47)	SB-11(5-5.5)	SB-11(29.5-30)	SB-12(12-14)
Detected Cor	npound Summary		Lab Sample Id	D4710-07	D4710-08	D4710-01	D4710-02	D4751-01	D4751-02	D4751-07
			Depth:	5 - 8 ft	28 - 30 ft	30 - 32 ft	45 - 47 ft	5 - 5.5 ft	29.5 - 30 ft	12 - 14 ft
			Source:	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH
			SDG:	D4710	D4710	D4710	D4710	D4751	D4751	D4751
		Unrestricted Use	Matrix:	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
		Soil Cleanup	Sampled:	10/26/2012 14:15	10/26/2012 14:55	10/25/2012 12:10	10/25/2012 12:40	11/6/2012 12:05	11/6/2012 12:10	11/5/2012 10:10
		Objectives	Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND		UNITS:							
	VOLATILES									
67-64-1	ACETONE	0.05	mg/kg	0.14	0.017 J	ND	ND	ND	ND	ND
71-43-2	BENZENE	0.06	mg/kg	0.028 J	0.0079	210	ND	ND	ND	ND
75-15-0	CARBON DISULFIDE		mg/kg	ND	ND	0.041	ND	0.0028 J	ND	ND
74-87-3	CHLOROMETHANE		mg/kg	ND	ND	ND	ND	ND	ND	ND
110-82-7	CYCLOHEXANE		mg/kg	ND	ND	0.057	ND	ND	ND	ND
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE		mg/kg	ND	ND	R	ND	ND	ND	ND
95-50-1	1,2-DICHLOROBENZENE	1.1	mg/kg	ND	ND	R	ND	ND	ND	ND
541-73-1	1,3-DICHLOROBENZENE	2.4	mg/kg	ND	ND	R	ND	ND	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND	R	ND	ND	ND	ND
75-71-8	DICHLORODIFLUOROMETHANE		mg/kg	ND	ND	ND	ND	ND	ND	ND
75-34-3	1,1-DICHLOROETHANE	0.27	mg/kg	ND	ND	ND	ND	ND	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	0.0031 J	0.002 J	540	ND	ND	ND	9.1
98-82-8	ISOPROPYLBENZENE (CUMENE)		mg/kg	ND	ND	29 J	ND	ND	ND	1
108-87-2	METHYLCYCLOHEXANE		mg/kg	ND	ND	0.18	ND	ND	ND	69
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	ND	ND	0.16 J	ND	ND	ND	ND
100-42-5	STYRENE		mg/kg	ND	ND	7.1 J	ND	ND	ND	0.29 J
79-34-5	1,1,2,2-TETRACHLOROETHANE		mg/kg	ND	ND	R	ND	ND	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	0.011 J	0.0059 J	510	ND	ND	ND	1.5
87-61-6	1,2,3-TRICHLOROBENZENE		mg/kg	ND	ND	R	ND	ND	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE		mg/kg	ND	ND	R	ND	ND	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	0.0021 J	0.0015 J	470	ND	ND	ND	11
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND	220	ND	ND	ND	5.7
	Total VOCs	NS	mg/kg	0.1842	0.0343	1986.538	ND	0.0028	ND	97.59
	SEMIVOLATILES		_							
92-52-4	BIPHENYL (DIPHENYL)		mg/kg	ND	ND	200	ND	ND	ND	10
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE		mg/kg	ND	ND	ND	ND	ND	ND	ND
86-74-8	CARBAZOLE		mg/kg	ND	ND	13	ND	ND	ND	ND
132-64-9	DIBENZOFURAN		mg/kg	ND	ND	43	ND	ND	ND	2.5
131-11-3	DIMETHYL PHTHALATE		mg/kg	0.161 J	ND	ND	ND	0.296 J	0.284 J	ND
84-74-2	DI-N-BUTYL PHTHALATE PAHs		mg/kg	ND	ND	ND	ND	ND	ND	ND
83-32-9	ACENAPHTHENE	20	mg/kg	ND	ND	320 J	ND	ND	ND	9.1
208-96-8	ACENAPHTHYLENE	100	mg/kg	ND	ND ND	370	0.2 J	ND	ND ND	50
120-12-7	ANTHRACENE	100	mg/kg	ND	ND	370	ND	ND	ND ND	22
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	0.29 J	ND ND	220	ND ND	ND	ND ND	11
50 55-5	DELIZO(LI)/IIII/ICENE	1	mg/Kg	0.27 3	ND	220	ND	ND	ND	11

C EL L	1 0''	1	Ir .: ID I	CD 0	CD 0	CD 10	CD 10	CD 11	CD 11	CD 10
Con Ed - Lud			Location ID:	SB- 9	SB- 9	SB-10	SB-10	SB-11	SB-11	SB-12
	l Analytical Data		Sample ID:	SB-9(5-8) D4710-07	SB-9(28-30)	SB-10(30-32)	SB-10(45-47)	SB-11(5-5.5)	SB-11(29.5-30)	SB-12(12-14)
Detected Con	npound Summary		Lab Sample Id		D4710-08	D4710-01	D4710-02	D4751-01	D4751-02	D4751-07
			Depth:	5 - 8 ft	28 - 30 ft	30 - 32 ft	45 - 47 ft	5 - 5.5 ft	29.5 - 30 ft	12 - 14 ft
			Source:	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH
			SDG:	D4710	D4710	D4710	D4710	D4751	D4751	D4751
		Unrestricted Use	Matrix:	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
		Soil Cleanup	Sampled:	10/26/2012 14:15	10/26/2012 14:55	10/25/2012 12:10	10/25/2012 12:40	11/6/2012 12:05	11/6/2012 12:10	11/5/2012 10:10
		Objectives	Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND		UNITS:							
	VOLATILES									
50-32-8	BENZO(A)PYRENE	1	mg/kg	0.28 J	ND	150 J	ND	ND	ND	7.2
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	0.38 J	ND	97 J	ND	ND	ND	5.5
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	0.17 J	ND	42 J	ND	ND	ND	2.1
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	ND	ND	32	ND	ND	ND	1.7 J
218-01-9	CHRYSENE	1	mg/kg	0.32 J	ND	220	ND	ND	ND	10
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	ND	ND	12	ND	ND	ND	ND
206-44-0	FLUORANTHENE	100	mg/kg	0.56	ND	400	ND	ND	ND	22
86-73-7	FLUORENE	30	mg/kg	ND	ND	550	ND	ND	ND	33
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	0.19 J	ND	37	ND	ND	ND	2
91-57-6	2-METHYLNAPHTHALENE		mg/kg	ND	ND	2800	ND	ND	ND	160
91-20-3	NAPHTHALENE	12	mg/kg	ND	ND	5700	0.21 J	ND	ND	310
85-01-8	PHENANTHRENE	100	mg/kg	0.32 J	ND	1500	ND	ND	ND	94
129-00-0	PYRENE	100	mg/kg	0.47	ND	640	ND	ND	ND	29
	Total PAHs		mg/kg	2.98	ND	13460	0.41	ND	ND	768.6
	Total SVOCs		mg/kg	3.141	ND	13716	0.41	0.296	0.284	781.1
	INORGANICS		Ŭ							
7429-90-5	ALUMINUM		mg/kg	7070	9490	3130	4490	4170 J	5890 J	4120 J
7440-38-2	ARSENIC	13	mg/kg	7.78	3	7.21	1.23	18.3 J	1.3 J	0.9 J
7440-39-3	BARIUM	350	mg/kg	157	96	30.9	55.6	156 J	63.3 J	31.9 J
7440-41-7	BERYLLIUM	7.2	mg/kg	0.11 J	0.39	0.04 J	0.2	ND	0.3	0.14 J
7440-43-9	CADMIUM	2.5	mg/kg	1.24	0.7	1.75	0.29	0.16 J	0.42	0.35
7440-70-2	CALCIUM		mg/kg	5480	2540	9540	13000	2260 J	12200 J	952 J
7440-47-3	CHROMIUM, TOTAL	30	mg/kg	16.7 J	18.4 J	5.38 J	8.9 J	9.93 J	11.7 J	9.3 J
7440-48-4	COBALT		mg/kg	7.34	11.9	3.4	5.38	6.31	6.51	6.88
7440-50-8	COPPER	50	mg/kg	44.1 J	20.5 J	26.5 J	11.8 J	69.6 J	12.1 J	11 J
7439-89-6	IRON		mg/kg	29900	28600	13200	12600	26800	14600	10300
7439-92-1	LEAD	63	mg/kg	462	18.1	215	8.81	30	10.5	45.3
7439-95-4	MAGNESIUM		mg/kg	2280	5050	5480	5450	1690 J	5740 J	1870 J
7439-96-5	MANGANESE	1600	mg/kg	284	414	277	290	803 J	319 J	191 J
7439-97-6	MERCURY	0.18	mg/kg	2.33	ND	0.381	ND	1.17	ND	ND
7440-02-0	NICKEL	30	mg/kg	21.2	31.1	12.2	14.3	17.5	16.6	11.2
7440-09-7	POTASSIUM		mg/kg	655	2730	352	1470	1500 J	2190 J	957 J
7782-49-2	SELENIUM	3.9	mg/kg	0.49	0.33 J	1.95	ND	1	ND	ND
7440-22-4	SILVER	2		0.79	0.67	0.15 J	0.09 J	0.51	0.14 J	0.1 J
7440-23-5	SODIUM			1150	148	4300	150	867 J	2520 J	5360 J
7440-28-0	THALLIUM			1.13	0.74 J	1.15	0.18 J	0.46 J	0.34 J	ND
7440-62-2	VANADIUM			16.7	22.5	7.59	13.1	14.8 J	15.3 J	12.7 J
7440-66-6	ZINC	109		268	50.8	506	22.9	88.4	27	18.6
57-12-5	CYANIDE	27	mg/kg	0.211 J	0.244 J	4.4	0.172 J	0.133 J	0.042 J	0.434
7782-49-2 7440-22-4 7440-23-5 7440-28-0 7440-62-2 7440-66-6	SELENIUM SILVER SODIUM THALLIUM VANADIUM ZINC	3.9 2 109	mg/kg mg/kg mg/kg mg/kg mg/kg mg/kg	0.49 0.79 1150 1.13 16.7 268	0.33 J 0.67 148 0.74 J 22.5 50.8	1.95 0.15 J 4300 1.15 7.59 506	ND 0.09 J 150 0.18 J 13.1 22.9	1 0.51 867 J 0.46 J 14.8 J 88.4	ND 0.14 J 2520 J 0.34 J 15.3 J 27	5.

Notes:

- (1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)
- (2) -- indicates no cleanup objective or background level is available.
- (3) ND indicates compound was not detected.
- (4) J indicates an estimated concentration.
- (5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.
- (6) NA indicates compound was not analyzed.

Dup of SB-14(45-47) SB-14 SB-14 Con Ed - Ludlow Site Location ID: SB-12 SB-13 SB-14 SB-16 SB-16 Sample ID: SB-12(36-38) SB-13(47-49) SB-14(38-40) SB-14(45-47) SB-14A(45-47) SB-16(30-33) SB-16(45-47) Validated Soil Analytical Data Lab Sample Id Detected Compound Summary D4751-08 D4751-09 D5300-03 D5300-06 D5300-07 D5300-01 D5300-02 47 - 49 ft 38 - 40 ft 45 - 47 ft 45 - 47 ft 30 - 33 ft 45 - 47 ft Depth: 36 - 38 ft CTECH Source: CTECH CTECH CTECH CTECH CTECH CTECH SDG: D4751 D5300 D4751 D5300 D5300 D5300 D5300 Unrestricted Use Matrix: SOIL SOIL SOIL SOIL SOIL SOIL SOIL Soil Cleanup Sampled: 11/5/2012 11:00 11/5/2012 15:05 12/21/2012 10:50 12/21/2012 11:50 12/21/2012 12:00 12/20/2012 16:10 12/20/2012 17:10 Objectives Validated: 1/23/2013 1/23/2013 1/23/2013 1/23/2013 1/23/2013 1/23/2013 1/23/2013 COMPOUND UNITS: CAS NO. VOLATILES 67-64-1 ACETONE 0.05 ND ND ND 0.027 J 0.026 J ND 0.071 mg/kg 71-43-2 BENZENE 0.06 ND 0.085 ND ND ND 17.2 ND mg/kg 75-15-0 CARBON DISULFIDE mg/kg ND 0.0026 J ND ND ND ND ND 74-87-3 CHLOROMETHANE --mg/kg ND ND ND ND ND 0.32 JND 110-82-7 CYCLOHEXANE ND 0.0017 J ND ND ND ND ND mg/kg 96-12-8 1.2-DIBROMO-3-CHLOROPROPANE mg/kg ND ND ND ND R 95-50-1 1,2-DICHLOROBENZENE 1.1 mg/kg ND ND ND ND ND ND ND 541-73-1 1,3-DICHLOROBENZENE 2.4 ND ND ND ND mg/kg ND ND ND 106-46-7 1,4-DICHLOROBENZENE 1.8 ND ND ND ND ND ND ND mg/kg 75-71-8 DICHLORODIFLUOROMETHANE ND ND ND ND ND ND ND mg/kg 75-34-3 0.27 1.1-DICHLOROETHANE ND ND ND ND ND ND ND mg/kg 100-41-4 ETHYLBENZENE 0.46 J 43.9 0.0045 J mg/kg ND ND ND ND 98-82-8 ISOPROPYLBENZENE (CUMENE) ND 0.045 ND ND ND ND 3.3 mg/kg 108-87-2 METHYLCYCLOHEXANE 0.0071 0.18 J mg/kg ND ND ND ND ND 75-09-2 METHYLENE CHLORIDE 0.05 mg/kg ND ND ND ND ND ND ND 100-42-5 0.0038 J STYRENE mg/kg ND ND ND ND 8.1 ND 79-34-5 1,1,2,2-TETRACHLOROETHANE ND ND ND ND ND ND mg/kg ND 108-88-3 TOLUENE 0.7 ND 0.053 ND ND ND 46.5 ND mg/kg 1,2,3-TRICHLOROBENZENE 87-61-6 mg/kg ND ND ND ND ND ND ND 120-82-1 1,2,4-TRICHLOROBENZENE ND ND ND ND ND ND ND mg/kg XYLMP M,P-XYLENE (SUM OF ISOMERS) 0.26 38.4 0.0039 J mg/kg ND 0.23 ND ND ND 95-47-6 O-XYLENE (1.2-DIMETHYLBENZENE) 0.26 ND ND 18.2 0.0022 J mg/kg ND 0.16 ND Total VOCs NS mg/kg ND 1.0482 ND 0.027 0.026 176.1 0.0816 SEMIVOLATILES 92-52-4 BIPHENYL (DIPHENYL) ND ND ND ND ND 1.6 ND mg/kg 117-81-7 BIS(2-ETHYLHEXYL) PHTHALATE mg/kg ND ND ND ND ND ND ND 86-74-8 CARBAZOLE ND ND ND ND ND ND ND mg/kg 132-64-9 DIBENZOFURAN mg/kg ND ND ND ND ND 0.53 ND 131-11-3 DIMETHYL PHTHALATE 0.455 0.403 0.57 0.32 J 0.34 J 0.25 J 0.4 J mg/kg 84-74-2 DI-N-BUTYL PHTHALATE ND ND ND ND ND ND ND mg/kg PAHs 83-32-9 20 ACENAPHTHENE mg/kg ND 0.32 J ND ND ND 1.8 ND 208-96-8 ACENAPHTHYLENE 100 ND 0.24 J ND ND ND 4.7 ND mg/kg 120-12-7 ANTHRACENE 100 ND 0.31 J ND ND ND 2.5 ND mg/kg

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mg/kg

ND

56-55-3

BENZO(A)ANTHRACENE

0.26 J

ND

ND

ND

ND

								Dup of SB-14(45-47)	1	
Con Ed - Lud	low Site		Location ID:	SB-12	SB-13	SB-14	SB-14	SB-14	SB-16	SB-16
	l Analytical Data		Sample ID:	SB-12(36-38)	SB-13(47-49)	SB-14(38-40)	SB-14(45-47)	SB-14A(45-47)	SB-16(30-33)	SB-16(45-47)
	npound Summary		Lab Sample Id	D4751-08	D4751-09	D5300-03	D5300-06	D5300-07	D5300-01	D5300-02
	1		Depth:	36 - 38 ft	47 - 49 ft	38 - 40 ft	45 - 47 ft	45 - 47 ft	30 - 33 ft	45 - 47 ft
			Source:	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH	CTECH
			SDG:	D4751	D4751	D5300	D5300	D5300	D5300	D5300
		Unrestricted Use	Matrix:	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
		Soil Cleanup	Sampled:	11/5/2012 11:00	11/5/2012 15:05	12/21/2012 10:50	12/21/2012 11:50	12/21/2012 12:00	12/20/2012 16:10	12/20/2012 17:10
		Objectives	Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	Objectives	UNITS:	1/20/2010	1/25/2015	1/23/2013	1/20/2010	1/23/2013	1/25/2015	1/23/2013
	VOLATILES									
50-32-8	BENZO(A)PYRENE	1 1	mg/kg	ND	0.2 J	ND	ND	ND	1.2	ND
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	ND	0.16 J	ND	ND	ND	0.94	ND
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	ND	ND	ND	ND	ND	0.31 J	ND
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	ND	ND	ND	ND	ND	0.38 J	ND
218-01-9	CHRYSENE	1	mg/kg	ND	0.26 J	ND	ND	ND	1.6	ND
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	ND	ND	ND	ND	ND	ND	ND
206-44-0	FLUORANTHENE	100	mg/kg	ND	0.49	ND	ND	ND	2.5	ND
86-73-7	FLUORENE	30	mg/kg	ND	0.32 J	ND	ND	ND	4.5	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	ND	ND	ND	ND	ND	0.2 J	ND
91-57-6	2-METHYLNAPHTHALENE		mg/kg	ND	1.3	ND	ND	ND	14.9	ND
91-20-3	NAPHTHALENE	12	mg/kg	ND	2.8	ND	ND	ND	33.7	ND
85-01-8	PHENANTHRENE	100	mg/kg	ND	1.3	ND ND	ND	ND	9.9	ND
129-00-0	PYRENE	100	mg/kg	ND	0.62	ND	ND	ND	3.3	ND
123-00-0	TRENE	100	mg/kg	ND	0.02	ND	ND	ND	3.3	ND
	Total PAHs		mg/kg	ND	8.58	ND	ND	ND	84.03	ND
	Total SVOCs		mg/kg	0.455	8.983	0.57	0.32	0.34	86.41	0.4
	INORGANICS									
7429-90-5	ALUMINUM		mg/kg	2800 J	9270 J	7400 J	2940 J	3060 J	14400 J	6040 J
7440-38-2	ARSENIC	13	mg/kg	0.49 J	1.96 J	0.97 J	0.71 J	0.89 J	5.98	1.18
7440-39-3	BARIUM	350	mg/kg	19.1 J	124 J	54.1 J	23.7 J	26.6 J	83.1 J	74.4 J
7440-41-7	BERYLLIUM	7.2	mg/kg	0.04 J	0.47	0.39	0.14 J	0.14 J	0.32 J	0.33
7440-43-9	CADMIUM	2.5	mg/kg	0.18	0.71	ND	ND	ND	ND	ND
7440-70-2	CALCIUM		mg/kg	10100 J	19200 J	2180 J	9830 J	9770 J	24700 J	12500 J
7440-47-3	CHROMIUM, TOTAL	30	mg/kg	7.04 J	19.9 J	16.3 J	ND	ND	12.4 J	12.9 J
7440-48-4	COBALT		mg/kg	3.39	7.91	7.86	3.26	3.53	3.36	7.11
7440-50-8	COPPER	50	mg/kg	9.25 J	18.3 J	7.99	5.12	5.88	5.21	9.59
7439-89-6	IRON		mg/kg	6990	18600	18000 J	8120 J	8770 J	15400 J	15800 J
7439-92-1	LEAD	63	mg/kg	3.88	15.2	10.1	4.47	4.63	25.8	8.35
7439-95-4	MAGNESIUM		mg/kg	4920 J	3760 J	3730 J	4070 J	3950 J	13500 J	5840 J
7439-96-5	MANGANESE	1600	mg/kg	209 J	1240 J	424 J	290 Ј	296 J	652 J	443 J
7439-97-6	MERCURY	0.18	mg/kg	ND	0.02	0.005 J	ND	0.003 J	0.003 J	0.005 J
7440-02-0	NICKEL	30	mg/kg	11.5	21.4	18.7	8.25	8.93	10.1	19.1
7440-09-7	POTASSIUM		mg/kg	581 J	2340 Ј	2720 J	832 J	824 J	1040 J	2050 J
7782-49-2	SELENIUM	3.9	mg/kg	ND	1.06	0.76 J	ND	ND	0.49 J	0.63 J
7440-22-4	SILVER	2	mg/kg	ND	0.34	0.51 J	0.17 J	0.15 J	0.23 J	0.37 J
7440-23-5	SODIUM		mg/kg	1220 J	1280 J	332	101	97.1 J	6400	1100
7440-28-0	THALLIUM		mg/kg	0.13 J	0.81 J	ND	ND	ND	ND	ND
7440-62-2	VANADIUM		mg/kg	9.49 J	18.3 J	19.3	8.87	10.2	19.1	16.6
7440-66-6	ZINC	109	mg/kg	14.9	33.7	26.9 J	13 J	12.9 J	12.6 J	25.1 J
57-12-5	CYANIDE	27	mg/kg	0.069 J	0.093 J	0.241 J	0.219 J	0.299	0.909	0.294 J

Notes:

- (1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)
- (2) -- indicates no cleanup objective or background level is available.
- (3) ND indicates compound was not detected.
- (4) J indicates an estimated concentration.
- (5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.
- (6) NA indicates compound was not analyzed.

Table 3
Summary of Soil Analytical Data
Former Ludlow Street Works Site
Consolidated Edison Company of New York
Remedial Investigation - October 2012 through April 2013

Con Ed - Luc	llow Site		Location ID:	SB-17	SB-17	TP-6FLOOR
Validated So	il Analytical Data		Sample ID:	SB-17(33-35)	SB-17(45-47)	TP-6FLOOR
Detected Cor	npound Summary		Lab Sample Id	D5300-08	D5300-09	D4751-10
	1		Depth:	33 - 35 ft	45 - 47 ft	-
			Source:	CTECH	CTECH	CTECH
			SDG:	D5300	D5300	D4751
		Unrestricted Use	Matrix:	SOIL	SOIL	SOIL
		Soil Cleanup	Sampled:	12/21/2012 15:00	12/21/2012 15:30	11/7/2012 10:30
		Objectives	Validated:	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	·	UNITS:			
	VOLATILES					
67-64-1	ACETONE	0.05	mg/kg	ND	0.011 J	0.026 J
71-43-2	BENZENE	0.06	mg/kg	ND	ND	ND
75-15-0	CARBON DISULFIDE		mg/kg	ND	ND	ND
74-87-3	CHLOROMETHANE		mg/kg	ND	ND	ND
110-82-7	CYCLOHEXANE		mg/kg	ND	ND	ND
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE		mg/kg	R	ND	R
95-50-1	1,2-DICHLOROBENZENE	1.1	mg/kg	ND	ND	R
541-73-1	1,3-DICHLOROBENZENE	2.4	mg/kg	ND	ND	R
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND	R
75-71-8	DICHLORODIFLUOROMETHANE		mg/kg	ND	ND	ND
75-34-3	1,1-DICHLOROETHANE	0.27	mg/kg	ND	ND	0.0066 J
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)		mg/kg	ND	ND	R
108-87-2	METHYLCYCLOHEXANE		mg/kg	ND	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	ND	ND	ND
100-42-5	STYRENE		mg/kg	ND	ND	ND
79-34-5	1,1,2,2-TETRACHLOROETHANE		mg/kg	ND	ND	R
108-88-3	TOLUENE	0.7	mg/kg	ND	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE		mg/kg	ND	ND	R
120-82-1	1,2,4-TRICHLOROBENZENE		mg/kg	ND	ND	R
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND	ND
	Total VOCs	NS	mg/kg	ND	0.011	0.0326
02.52.4	SEMIVOLATILES DEPARTMENT OF THE PROPERTY OF T			N.T.		
92-52-4	BIPHENYL (DIPHENYL)		mg/kg	ND	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE		mg/kg	ND	0.48	ND
86-74-8	CARBAZOLE		mg/kg	ND	ND	ND ND
132-64-9	DIBENZOFURAN		mg/kg	ND	ND	
131-11-3	DIMETHYL PHTHALATE		mg/kg	0.35 ND	0.39	0.632 ND
84-74-2	DI-N-BUTYL PHTHALATE PAHs		mg/kg	ND	0.15 J	ND
83-32-9	ACENAPHTHENE	20	mg/kg	ND	ND	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	ND ND	ND ND	ND ND
120-12-7	ANTHRACENE	100	mg/kg	ND ND	ND	ND ND
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	ND ND	ND ND	0.35 J

C F1 I I	1 00	Т	ly .: TD	CD 17	CD 17	TD (FLOOD
Con Ed - Lud			Location ID:	SB-17	SB-17	TP-6FLOOR
	l Analytical Data		Sample ID: Lab Sample Id	SB-17(33-35) D5300-08	SB-17(45-47) D5300-09	TP-6FLOOR D4751-10
Detected Con	npound Summary				45 - 47 ft	D4/31-10
			Depth: Source:	33 - 35 ft CTECH	CTECH	CTECH
			Source: SDG:	D5300	D5300	D4751
		Unrestricted Use	Matrix:	SOIL	SOIL	SOIL
		Soil Cleanup	Sampled:	12/21/2012 15:00	12/21/2012 15:30	11/7/2012 10:30
		Objectives	Validated:	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	Objectives	UNITS:	1/23/2013	1/23/2013	1/23/2013
CAS NO.	VOLATILES		UNITS.			
50-32-8	BENZO(A)PYRENE	1	mg/kg	ND	ND	0.41 J
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	ND ND	ND ND	0.41 3
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	ND ND	ND ND	0.29 J
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	ND ND	ND ND	ND
218-01-9	CHRYSENE	1	mg/kg	ND ND	ND ND	0.4 J
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	ND ND	ND ND	ND
206-44-0	FLUORANTHENE	100	mg/kg	ND ND	ND ND	0.75
86-73-7	FLUORENE	30	mg/kg	ND	ND	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	ND ND	ND ND	0.27 J
91-57-6	2-METHYLNAPHTHALENE		mg/kg	ND	0.19 J	ND
91-20-3	NAPHTHALENE	12	mg/kg	ND ND	0.15 J	ND ND
85-01-8	PHENANTHRENE	100	mg/kg	ND	ND	0.45 J
129-00-0	PYRENE	100	mg/kg	ND ND	ND ND	0.69
127 00 0	TREAT	100	mg/kg	ND	ND	0.07
	Total PAHs		mg/kg	ND	0.44	4.11
	Total SVOCs		mg/kg	0.35	1.46	4.742
	INORGANICS					
7429-90-5	ALUMINUM		mg/kg	2860 J	9120 J	9100 J
7440-38-2	ARSENIC	13	mg/kg	0.68 J	0.73 J	14.5 J
7440-39-3	BARIUM	350	mg/kg	33.3 J	50.8 J	122 J
7440-41-7	BERYLLIUM	7.2	mg/kg	0.1 J	0.06 J	ND
7440-43-9	CADMIUM	2.5	mg/kg	ND	ND	1.73
7440-70-2	CALCIUM		mg/kg	1980 J	5560 J	6580 J
7440-47-3	CHROMIUM, TOTAL	30	mg/kg	ND	25.5 J	28.7 J
7440-48-4	COBALT		mg/kg	4.01	6.27	9.95
7440-50-8	COPPER	50	mg/kg	8.81	31.1	173 J
7439-89-6	IRON		mg/kg	7530 J	13000 J	39400
7439-92-1	LEAD	63	mg/kg	2.92	20.3	356
7439-95-4	MAGNESIUM		mg/kg	2040 J	3950 J	4080 J
7439-96-5	MANGANESE	1600	mg/kg	519 J	332 J	345 J
7439-97-6	MERCURY	0.18	mg/kg	0.003 J	0.014	6.2
7440-02-0	NICKEL	30	mg/kg	11.8	17.6	29.2
7440-09-7	POTASSIUM		mg/kg	730 J	1420 J	901 J
7782-49-2	SELENIUM	3.9	mg/kg	0.38 J	0.51 J	2.18
7440-22-4	SILVER	2	mg/kg	0.17 J	0.18 J	0.39
7440-23-5	SODIUM		mg/kg	188	913	2090 J
7440-28-0	THALLIUM		mg/kg	ND	ND	1.46
7440-62-2	VANADIUM		mg/kg	10.3	22.4	26 J
7440-66-6	ZINC	109	mg/kg	13.1 J	28.6 J	315
57-12-5	CYANIDE	27	mg/kg	0.172 J	0.257 J	0.297 J

Notes:

- (1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)
- (2) -- indicates no cleanup objective or background level is available.
- (3) ND indicates compound was not detected.
- (4) J indicates an estimated concentration.
- (5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.
- (6) NA indicates compound was not analyzed.

Remedial Investigation - October 2012 through April 2013

C 11.1 . 1	1.0.1		lr .: rp l) (IV. 1	3.037.1	3.037.0
Consolidated			Location ID:	MW-1	MW-1	MW-2
Ludlow Stree			Sample ID:	MW-1	MW-1-20130405	MW-2-20130408
	oundwater Analytical Data		Lab Sample Id:	D4947-01	E1768-02	E1768-07
Detected Cor	mpound Summary		Source:	CTECH	CTECH	CTECH
		NYSDEC	SDG:	D4947	E1768	E1768
		Class GA	Matrix:	WATER	WATER	WATER
		Groundwater	Sampled:	11/26/2012	4/5/2013	4/8/2013
		Standards/Guidance	Validated:	1/23/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	Values (1)	UNITS:			
	VOLATILES					
67-64-1	ACETONE	5	ug/l	ND	7.8 J	8.7 J
74-83-9	BROMOMETHANE	5	ug/l	ND	ND	ND
110-82-7	CYCLOHEXANE		ug/l	ND	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	5	ug/l	ND	ND	ND
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	50 (G)	ug/l	ND	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	30 (G) 	ug/l	ND ND	ND ND	ND
108-10-1			-			
	METHYLCYCLOHEXANE	 5	ug/l	ND ND	ND ND	ND ND
100-42-5	STYRENE	5	ug/l	ND	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	10	ug/l	ND	ND	ND
	BTEX					. —
71-43-2	BENZENE	1	ug/l	ND	ND	ND
100-41-4	ETHYLBENZENE	5	ug/l	ND	ND	ND
108-88-3	TOLUENE	5	ug/l	ND	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	5	ug/l	ND	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	5	ug/l	ND	ND	ND
	CVOCs					
74-87-3	CHLOROMETHANE	5	ug/l	ND	ND	ND
156-59-2	CIS-1,2-DICHLOROETHYLENE	5	ug/l	ND	ND	0.49 J
127-18-4	TETRACHLOROETHYLENE(PCE)	5	ug/l	ND	ND	ND
,, -			1.8.1			
	Total VOCs			ND	7.8	9.19
	SEMIVOLATILES					
98-86-2	ACETOPHENONE		ug/l	ND	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	5	ug/l	ND	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	5	ug/l	ND	ND	ND
86-74-8	CARBAZOLE		ug/l	ND	ND	ND
132-64-9	DIBENZOFURAN		ug/l	ND	ND	ND
91-94-1	3,3'-DICHLOROBENZIDINE	5	ug/l	ND	ND	ND
99-09-2	3-NITROANILINE	5	ug/l	ND	ND	ND
108-95-2	PHENOL	1	ug/l	ND	ND	ND
	PAHs	-	ug/1	1,2	1,2	1,12
83-32-9	ACENAPHTHENE	20 (G)	ug/l	ND	ND	ND
208-96-8	ACENAPHTHYLENE	20 (G) 	ug/l	ND ND	ND ND	ND ND
120-12-7	ANTHRACENE	50 (G)		ND ND	ND ND	ND ND
56-55-3			ug/l			
	BENZO(A)ANTHRACENE	0.002 (G)	ug/l	ND	ND	ND
50-32-8	BENZO(A)PYRENE	ND	ug/l	ND	ND	ND
205-99-2	BENZO(B)FLUORANTHENE	0.002 (G)	ug/l	ND	ND	ND
218-01-9	CHRYSENE	0.002 (G)	ug/l	ND	ND	ND
	FLUORANTHENE	50 (G)	ug/l	ND	ND	ND
	FLUORENE	50 (G)	ug/l	ND	ND	ND
91-57-6	2-METHYLNAPHTHALENE		ug/l	ND	ND	ND
91-20-3	NAPHTHALENE	10 (G)	ug/l	ND	ND	ND
85-01-8	PHENANTHRENE	50 (G)	ug/l	ND	ND	ND
	PYRENE	50 (G)	ug/l	ND	ND	ND
129-00-0	TREAL	` '	1			
	Total PAHs	, ,		ND	ND	ND

Indicates concentration exceeds standard or guidance value.

(G) Indicates guidance value.

No standard or guidance value available. Indicates compound was not detected. NS ND Indicates an estimated concentration.

Micrograms per liter ug/L

Remedial Investigation - October 2012 through April 2013

Consolidated	1 Edison		Location ID:	MW-1	MW-1	MW-2
Ludlow Stree			Sample ID:	MW-1	MW-1-20130405	MW-2-20130408
	roundwater Analytical Data		Lab Sample Id:	D4947-01	E1768-02	E1768-07
	mpound Summary		Source:	CTECH	CTECH	CTECH
Detected Co.	mpound summary	NYSDEC	SDG:	D4947	E1768	E1768
		Class GA	Matrix:	WATER	WATER	WATER
		Groundwater	Sampled:	11/26/2012	4/5/2013	4/8/2013
			Validated:	1/23/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	Values (1)	UNITS:	1/20/2010	170072010	110012010
CAS NO.	VOLATILES	values	UNIIS.			
	INORGANICS	 				
7429-90-5	ALUMINUM		ug/l	331 J	2380 Ј	368 J
7440-36-0	ANTIMONY	3	ug/l	4.98 J	ND	9.48 J
7440-38-2	ARSENIC	25	ug/l	6.1	7.8 J	ND
7440-39-3	BARIUM	1000	ug/l	159 J	129	1990
7440-41-7	BERYLLIUM	3 (G)	ug/l	ND	ND	ND
7440-43-9	CADMIUM	5	ug/l	ND ND	R	1.22 J
7440-70-2	CALCIUM	3	ug/l	91200 J	59000 J	262000 J
7440-70-2	CHROMIUM, TOTAL	50	ug/l ug/l	91200 J ND	37.6 J	202000 J ND
7440-47-3	COBALT	30	ug/l ug/l	ND ND	37.6 J ND	6.91 J
7440-50-8	COPPER	200	-	6.08 J	7.59 J	27.4 J
7439-89-6	IRON	300	ug/l	5890 J	8770	6860
7439-89-0	LEAD	25	ug/l	1.34 J	6.36 J	10.1 J
			ug/l			
7439-95-4 7439-96-5	MAGNESIUM MANGANESE	35000 (G)	ug/l	22100 J	13400 J	33200 J
7439-96-3		300	ug/l	4090 J	2200 J	2160 J
	MERCURY	0.7	ug/l	0.092 J	ND	ND
7440-02-0	NICKEL POTE SCHIP	100	ug/l	ND	20.9 J	ND
7440-09-7	POTASSIUM		ug/l	8980 J	6390	113000
7782-49-2	SELENIUM	10	ug/l	ND	ND	ND
7440-22-4	SILVER	50	ug/l	ND	ND	ND
7440-23-5	SODIUM	20000	ug/l	361000 J	272000 J	39700000 J
7440-28-0	THALLIUM	0.5 (G)	ug/l	ND	ND	ND
7440-62-2	VANADIUM		ug/l	ND	ND	ND
7440-66-6	ZINC	2000 (G)	ug/l	ND	18.4 J	35.4 J
57-12-5	CYANIDE	200	ug/L	ND	ND	224
7420 00 5	DISSOLVED METALS	ļ	σ.		ND	
7429-90-5	ALUMINUM		ug/L		ND	
7440-36-0	ANTIMONY	3	ug/L		ND	
7440-38-2	ARSENIC	25	ug/L		ND	
7440-39-3	BARIUM	1000	ug/L		64.9	
7440-41-7	BERYLLIUM	3 (G)	ug/L		ND	
7440-43-9	CADMIUM	5	ug/L		R	
7440-70-2	CALCIUM		ug/L		54500 J	
7440-47-3	CHROMIUM	50	ug/L		ND	
7440-48-4	COBALT		ug/L		ND	
7440-50-8	COPPER	200	ug/L		ND	
7439-89-6	IRON	300	ug/L		ND	
7439-92-1	LEAD	25	ug/L		ND	
7439-95-4	MAGNESIUM	35000 (G)	ug/L		10700 J	
7439-96-5	MANGANESE	300	ug/L		680 J	
7439-97-6	MERCURY	0.7	ug/L		ND	
7440-02-0	NICKEL	100	ug/L		ND	
7440-09-7	POTASSIUM		ug/L		6360	
7782-49-2	SELENIUM	10	ug/L		ND	
7440-23-5	SODIUM	20000	ug/L		271000 J	
7440-62-2	VANADIUM		ug/L		ND	
7440-66-6	ZINC	2000 (G)	ug/L		9.09 J	

Notes:

Indicates concentration exceeds standard or guidance value.

(G)

Indicates guidance value.

No standard or guidance value available. NS Indicates compound was not detected. Indicates an estimated concentration. ND

ug/L Micrograms per liter

Remedial Investigation - October 2012 through April 2013

Ladiow Street Size	Consolidated	1 Edison		Location ID:	MW-3	MW-3	MW-3A
Lab Sample Is D4947-49DI. E1768-08 E1768-09 E							
Detected Compound Summary							
NYSDEC Class GA Groundwater Class GA Groundwater Standard/Stofindance Values Groundwater Standard/Stofindance Values Groundwater Standard/Stofindance Values Groundwater Mystandance M		•					
Class GA Groundwater Standards/Guidance WATER WATER WATER Standards/Guidance Validance	Detected Co	inpound Summary	NYSDEC				
CAS NO. COMPOUND Values							
Standard/Sciulance Validated: 1/23/2013 4/30/201							
CAS NO. COMPOUND Values							
VOLATILES STACETONE STAC	CAS NO	COMPOLIND		1	1/23/2013	4/30/2013	4/30/2013
10.74 ACETONE	CAS NO.		v aiues	UNIIS.			
Table Tabl	67-64-1		5	110/1	ND	ND	ND
11082-7 CYCLOHEXANE							
98-82-8 ISOPROPYLEENZENE (CUMENE) 5 ug/l 7 3.9 J 4.6 J 7 3.9 J 4.0 J 7 3.9 J 4.0 J 7 3.9 J 4.2 J 7 3.9 J 4.2 J 7 3.9 J 4.0 J 4.2 J 7 3.9 J 4.2 J 7 3.9 J 4.0 J 4.2 J 7 3.9 J 4.0 J 4.2 J 7 3.9 J 4.0 J 4							
Transport Tran				-			
108-10-1 METHYL ISOBUTYL RETONE ug/l 10 ND ND ND ND 108-87-2 METHYLCYCLOHEXANE ug/l 59 41.5 42.6 16340-44 TERT-BUTYL METHYL ETHER 10 ug/l ND 1.7 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 1.8 3 3 1.8 3 3 3 3 3 3 3 3 3			_	-			
108-87-2		` '	` ′				
100-42-5 STYRENE 5 ug/l 59 41.5 42.6 1634-044 TERT-BUTYL METHYL ETHER 10 ug/l ND 1.7 J 1.8 J TERT 1.8 J TERT 1.5 ug/l 350 360 460 460 180 100-41-4							
1634-04-4 TERT-BUTYL METHYL ETHER 10				-			
BTEX				-			
1	1001017		10	ug/1	110	1./ 3	1.0 3
100-14-14	71_43_2		1	110/1	350	360	460
108-88-3 TOLUENE 5 ug/l 280 260 310 310 371 371 372 372 373 373 373 373 373 374							
XYLENE M.P. XYLENE (SUM OF ISOMERS) 5 ug/l 190 160 180 180 160 180 180 160 180 180 160 180 180 160 180 180 160 180 180 160 180 180 160 180 180 160 180 180 160 180 1				-			
190 160 180							
CVOCs				-			
Total VOCs	75-47-0		3	ug/1	170	100	100
156-59-2 CIS-1,2-DICHLOROETHYLENE 5 ug/l 0.45 J ND ND ND	74-87-3		5	110/1	ND	ND	ND
127-18-4 TETRACHLOROETHYLENE(PCE) 5 ug/l 0.42 J ND ND							
Total VOCs				-			
SEMIVOLATILES	127 10 1	TETRICIENDOROETHT EEL (E(T CE)	5	ug/1	0.12 3	TVD	TVD
See Sec 2 ACETOPHENONE					1562.37	1291.7	1555.3
92-52-4 BIPHENYL (DIPHENYL) 5 ug/l 25.1 J 25.2 36.2							
117-81-7 BIS(2-ETHYLHEXYL) PHTHALATE 5 ug/l ND ND ND							
132-64-9 DIBENZOFURAN ug/l 4.5 J 6.1 J 8.9 J 194-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND 108-95-2 PHENOL 1 ug/l ND ND ND 208-96-8 ACENAPHTHENE 5 ug/l ND ND ND 208-96-8 ACENAPHTHYLENE 5 ug/l ND 12.1 205-99-2 BENZO(A)ANTHRACENE 50 (G) ug/l ND ND 12.1 205-99-2 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND 8.5 J 218-01-9 CHRYSENE 0.002 (G) ug/l ND 6.4 J 219-37-7 FLUORENE 50 (G) ug/l 7.6 J 217-8 7.6 J 17.8 219-90-0 PYRENE 50 (G) ug/l 34.7 J 3030.5 3255.7 4985.6 Total PAHs 3030.5 3255.7 4985.6		, , , , , , , , , , , , , , , , , , , ,					
132-64-9 DIBENZOFURAN ug/l 4.5 J 6.1 J 8.9 J 91-94-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND 99-09-2 3-NITROANILINE 5 ug/l ND ND ND PHENOL 1 ug/l ND ND ND PAHS		· · · · · · · · · · · · · · · · · · ·	5				
91-94-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND ND ND ND ND N							
99-09-2 3-NITROANILINE 5 ug/l ND				-			
1				-			
PAHS				-			
Sa-32-9 ACENAPHTHENE 20 (G) ug/l 17.1 J 22.5 32.2	108-95-2		1	ug/l	ND	ND	ND
208-96-8 ACENAPHTHYLENE ug/l 84.2 75.6 J 130 J 120-12-7 ANTHRACENE 50 (G) ug/l 10.6 J 21.9 27.4 56-55-3 BENZO(A)ANTHRACENE 0.002 (G) ug/l ND 12.1 12.2 50-32-8 BENZO(A)PYRENE ND ug/l ND 8.5 J 8.5 J 205-99-2 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND 6.4 J 6.5 J 218-01-9 CHRYSENE 0.002 (G) ug/l ND 12.3 12.2 206-44-0 FLUORANTHENE 50 (G) ug/l 7.6 J 17.8 19.9 86-73-7 FLUORENE 50 (G) ug/l 34.7 J 47.3 65.1 91-57-6 2-METHYLNAPHTHALENE ug/l 320 630 630 91-20-3 NAPHTHALENE 10 (G) ug/l 2500 2300 3900 85-01-8 PHENANTHRENE 50 (G) ug/l 46.2 J 72.9 110 129-00-0 PYRENE 50 (G) ug/l 10.1 J 28.4 31.6 Total PAHs 3030.5 3255.7 4985.6							
120-12-7 ANTHRACENE 50 (G) ug/l 10.6 J 21.9 27.4			20 (G)	-			
56-55-3 BENZO(A)ANTHRACENE 0.002 (G) ug/l ND 12.1 12.2 50-32-8 BENZO(A)PYRENE ND ug/l ND 8.5 J 8.5 J 205-99-2 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND 6.4 J 6.5 J 218-01-9 CHRYSENE 0.002 (G) ug/l ND 12.3 12.2 206-44-0 FLUORANTHENE 50 (G) ug/l 7.6 J 17.8 19.9 86-73-7 FLUORENE 50 (G) ug/l 34.7 J 47.3 65.1 91-57-6 2-METHYLNAPHTHALENE ug/l 320 630 630 91-20-3 NAPHTHALENE 10 (G) ug/l 2500 2300 3900 85-01-8 PHENANTHRENE 50 (G) ug/l 46.2 J 72.9 110 129-00-0 PYRENE 50 (G) ug/l 10.1 J 28.4 31.6 Total PAHs				-			
Decomposition Section			* *	-			
Description				-			
218-01-9 CHRYSENE 0.002 (G) ug/l ND 12.3 12.2		` /		-			
206-44-0 FLUORANTHENE 50 (G) ug/l 7.6 J 17.8 19.9				-			
86-73-7 FLUORENE 50 (G) ug/l 34.7 J 47.3 65.1 91-57-6 2-METHYLNAPHTHALENE ug/l 320 630 630 91-20-3 NAPHTHALENE 10 (G) ug/l 2500 2300 3900 85-01-8 PHENANTHRENE 50 (G) ug/l 46.2 J 72.9 110 129-00-0 PYRENE 50 (G) ug/l 10.1 J 28.4 31.6 Total PAHs			* *	-			
91-57-6 91-20-3 NAPHTHALENE 91-20-3 NAPHTHALEN				-			
91-20-3 NAPHTHALENE 85-01-8 PHENANTHRENE 129-00-0 PYRENE 10 (G) ug/l 2500 2300 3900 50 (G) ug/l 46.2 J 72.9 110 50 (G) ug/l 10.1 J 28.4 31.6 Total PAHs 3030.5 3255.7 4985.6			* *	-			
85-01-8 PHENANTHRENE							
129-00-0 PYRENE 50 (G) ug/l 10.1 J 28.4 31.6 Total PAHs 3030.5 3255.7 4985.6				-			
Total PAHs 3030.5 3255.7 4985.6				-			
	129-00-0	PYRENE	50 (G)	ug/l	10.1 J	28.4	31.6
Total SVOCs 3066 3 3297 1 5047 1		Total PAHs			3030.5	3255.7	4985.6
		Total SVOCs			3066.3	3297.1	5047.1

Notes:

Indicates concentration exceeds standard or guidance value.

(G) Indicates guidance value.

No standard or guidance value available. Indicates compound was not detected. NS ND Indicates an estimated concentration.

Micrograms per liter ug/L

Remedial Investigation - October 2012 through April 2013

Consolidated	1 Edison	I	Location ID:	MW-3	MW-3	MW-3A
Ludlow Stree			Sample ID:	MW-3		MW-3A-20130408
	oundwater Analytical Data		Lab Sample Id:	D4947-09DL	E1768-08	E1768-09
	mpound Summary		Source:	CTECH	CTECH	CTECH
Detected Co.	inpound Summary	NYSDEC	SDG:	D4947	E1768	E1768
		Class GA	Matrix:	WATER	WATER	WATER
		Groundwater	Sampled:	11/27/2012	4/8/2013	4/8/2013
	T		Validated:	1/23/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	Values (1)	UNITS:			
	VOLATILES					
	INORGANICS					
7429-90-5	ALUMINUM		ug/l	ND	96.1 J	81.4 J
7440-36-0	ANTIMONY	3	ug/l	5.31 J	ND	ND
7440-38-2	ARSENIC	25	ug/l	2.45 J	ND	4.25 J
7440-39-3	BARIUM	1000	ug/l	106 J	134	128
7440-41-7	BERYLLIUM	3 (G)	ug/l	ND	ND	ND
7440-43-9	CADMIUM	5	ug/l	0.32 J	0.52 J	0.5 R
7440-70-2	CALCIUM		ug/l	58900 J	81900 J	77000 J
7440-47-3	CHROMIUM, TOTAL	50	ug/l	ND	ND	ND
7440-48-4	COBALT		ug/l	ND	ND	ND
7440-50-8	COPPER	200	ug/l	7.44 J	7.59 J	6.54 J
7439-89-6	IRON	300	ug/l	1160 J	1600	1590
7439-92-1	LEAD	25	ug/l	6.72	40.3 J	37.6 J
7439-95-4	MAGNESIUM	35000 (G)	ug/l	21500 J	22600 J	21000 J
7439-96-5	MANGANESE	300	ug/l	1100 J	1500 J	1400 J
7439-97-6	MERCURY	0.7	ug/l	ND	ND	ND
7440-02-0	NICKEL	100	ug/l	ND	5.62 J	5.4 J
7440-09-7	POTASSIUM		ug/l	8370 J	15500	14300
7782-49-2	SELENIUM	10	ug/l	ND	ND	ND
7440-22-4	SILVER	50	ug/l	ND	ND	ND
7440-23-5	SODIUM	20000	ug/l	3480000 J	4620000 J	5420000 J
7440-28-0	THALLIUM	0.5 (G)	ug/l	ND	ND	ND
7440-62-2	VANADIUM		ug/l	ND	ND	ND
7440-66-6	ZINC	2000 (G)	ug/l	ND	20.8 J	21.1 J
57-12-5	CYANIDE	200	ug/L	123	36 J	166 J
	DISSOLVED METALS					
7429-90-5	ALUMINUM		ug/L			
7440-36-0	ANTIMONY	3	ug/L			
7440-38-2	ARSENIC	25	ug/L			
7440-39-3	BARIUM	1000	ug/L			
7440-41-7	BERYLLIUM	3 (G)	ug/L			
7440-43-9	CADMIUM	5	ug/L			
7440-70-2	CALCIUM		ug/L			
7440-47-3	CHROMIUM	50	ug/L			
7440-48-4	COBALT		ug/L			
7440-50-8	COPPER	200	ug/L			
7439-89-6	IRON	300	ug/L			
7439-92-1	LEAD	25	ug/L			
	MAGNESIUM	35000 (G)	ug/L			
	MANGANESE	300	ug/L			
7439-97-6	MERCURY	0.7	ug/L			
7440-02-0	NICKEL	100	ug/L			
7440-09-7	POTASSIUM		ug/L			
7782-49-2	SELENIUM	10	ug/L ug/L			
7440-23-5	SODIUM	20000	ug/L ug/L			
7440-23-3	VANADIUM	20000	ug/L ug/L			
7440-62-2	ZINC	2000 (G)	ug/L ug/L			
Notes:	ZITC	2000 (0)	ug/L		l	l

Notes:

Indicates concentration exceeds standard or guidance value.

(G)

Indicates guidance value.

No standard or guidance value available. NS Indicates compound was not detected. Indicates an estimated concentration. ND

ug/L Micrograms per liter

Remedial Investigation - October 2012 through April 2013

Consolidated	4 Edison		Location ID:	MW-4	MW-4	MW-5
Ludlow Stre			Sample ID:	MW-4 MW-4	MW-4-20130408	MW-5-20130405
	roundwater Analytical Data		Lab Sample Id:	D4947-07	E1768-06	E1768-03
	mpound Summary		Source:	CTECH	CTECH	CTECH
Detected Co	inpound Summary	NYSDEC	SDG:	D4947	E1768	E1768
		Class GA	Matrix:	WATER	WATER	WATER
		Groundwater	Sampled:			
		Standards/Guidance	Validated:	11/27/2012	4/8/2013	4/5/2013
CACNO	COMPOUND		1	1/23/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	Values (1)	UNITS:			
67.64.1	VOLATILES	5		ND	01.1.7	220. I
67-64-1	ACETONE	5	ug/l	ND	81.1 J	320 J
74-83-9	BROMOMETHANE	5	ug/l	ND	ND	ND
110-82-7	CYCLOHEXANE		ug/l	ND	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	5	ug/l	ND	ND	1.8 J
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	50 (G)	ug/l	ND	ND	29.2 J
108-10-1	METHYL ISOBUTYL KETONE		ug/l	ND	ND	200 J
108-87-2	METHYLCYCLOHEXANE		ug/l	ND	ND	ND
100-42-5	STYRENE	5	ug/l	ND	ND	0.49 J
1634-04-4	TERT-BUTYL METHYL ETHER	10	ug/l	ND	ND	ND
	BTEX					
71-43-2	BENZENE	1	ug/l	ND	ND	35.8
100-41-4	ETHYLBENZENE	5	ug/l	ND	ND	12.1
108-88-3	TOLUENE	5	ug/l	ND	ND	36.9
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	5	ug/l	ND	ND	7.7 J
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	5	ug/l	ND	ND	8.1
	CVOCs					
74-87-3	CHLOROMETHANE	5	ug/l	ND	ND	ND
156-59-2	CIS-1,2-DICHLOROETHYLENE	5	ug/l	ND	ND	ND
127-18-4	TETRACHLOROETHYLENE(PCE)	5	ug/l	ND	ND	ND
	T I VOC			ND	01.1	(52.00
	Total VOCs SEMIVOLATILES			ND	81.1	652.09
98-86-2	ACETOPHENONE		a/l	ND	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	 5	ug/l ug/l	ND ND	ND ND	19.6
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	5	-	15.4 J	ND ND	ND
86-74-8	CARBAZOLE	3	ug/l	ND	ND ND	ND ND
132-64-9	DIBENZOFURAN		ug/l	ND ND	ND ND	ND ND
91-94-1		 5	ug/l			
	3,3'-DICHLOROBENZIDINE		ug/l	ND	ND	R
99-09-2	3-NITROANILINE	5	ug/l	ND	ND	R
108-95-2	PHENOL	1	ug/l	ND	ND	ND
83-32-9	PAHs ACENAPHTHENE	20 (C)	110/1	ND	ND	42.6
83-32-9 208-96-8		20 (G)	ug/l			
	ACENAPHTHYLENE		ug/l	ND	ND	18.1
120-12-7	ANTHRACENE DENZO(A) ANTHRACENE	50 (G)	ug/l	ND	ND	ND
56-55-3	BENZO(A) ANTHRACENE	0.002 (G)	ug/l	ND	ND	ND
50-32-8	BENZO(A)PYRENE	ND	ug/l	ND	ND	ND
205-99-2	BENZO(B)FLUORANTHENE	0.002 (G)	ug/l	ND	ND	ND
218-01-9	CHRYSENE	0.002 (G)	ug/l	ND	ND	ND
206-44-0	FLUORANTHENE	50 (G)	ug/l	ND	ND	ND
86-73-7	FLUORENE	50 (G)	ug/l	ND	ND	18
91-57-6	2-METHYLNAPHTHALENE		ug/l	ND	ND	42.1
91-20-3	NAPHTHALENE	10 (G)	ug/l	ND	ND	240 J
85-01-8	PHENANTHRENE	50 (G)	ug/l	ND	ND	21.2
129-00-0	PYRENE	50 (G)	ug/l	ND	ND	ND
	Total PAHs			ND	ND	382
	Total SVOCs			15.4	ND	401.6
Notes:						

Indicates concentration exceeds standard or guidance value.

(G) Indicates guidance value.

No standard or guidance value available. Indicates compound was not detected. NS ND Indicates an estimated concentration.

Micrograms per liter ug/L

Remedial Investigation - October 2012 through April 2013

Consolidated	LEdison		Location ID:	MW 4	MW-4	MW-5
				MW-4 MW-4	MW-4 MW-4-20130408	MW-5-20130405
Ludlow Street Site Validated Groundwater Analytical Data			Sample ID:		E1768-06	
	mpound Summary		Lab Sample Id: Source:	D4947-07 CTECH	CTECH	E1768-03 CTECH
Detected Col	mpound Summary	NYSDEC	Source: SDG:	D4947	E1768	E1768
		Class GA	Matrix:	WATER	WATER	WATER
		Groundwater	Sampled:	11/27/2012	4/8/2013	4/5/2013
			Validated:	1/23/2013	4/30/2013	4/30/2013
CACNO	COLEOTRIC	Values (1)	1	1/23/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	Values	UNITS:			
	VOLATILES INORGANICS					
7429-90-5	ALUMINUM		na/1	161000 J	32100 J	550 J
7440-36-0	ANTIMONY	3	ug/l	ND	ND	12.1 J
7440-38-2	ARSENIC	25	ug/l	206	40	4.4 J
7440-38-2	BARIUM	1000	ug/l	2170 J	523	916
	BERYLLIUM	3 (G)	ug/l	7.1 J	1.23 J	ND
7440-41-7	CADMIUM	5 (G) 5	ug/l	25	1.23 J 1.3 J	2.2 J
7440-43-9	CALCIUM		ug/l	67500 J	116000 J	2.2 J 267000 J
7440-70-2	CHROMIUM, TOTAL	50	ug/l	352 J	116000 J 113 J	267000 J ND
7440-47-3	COBALT	50 	ug/l	263	25.1 J	ND ND
7440-48-4	COPPER	200	ug/l	822 J	25.1 J 205 J	ND 5.86 J
7439-89-6			ug/l			
	IRON	300	ug/l	520000 J	55300	4550
7439-92-1	LEAD	25	ug/l	10700	1200 J	232 J
7439-95-4	MAGNESIUM	35000 (G)	ug/l	53000 J	24500 J	59800 J
7439-96-5	MANGANESE	300	ug/l	6820 J	1210 J	1580 J
7439-97-6	MERCURY	0.7	ug/l	16.85	20.182	ND
7440-02-0	NICKEL	100	ug/l	689	112 J	ND
7440-09-7	POTASSIUM		ug/l	23300 J	11800	93500
7782-49-2	SELENIUM	10	ug/l	26	16.1	ND
7440-22-4	SILVER	50	ug/l	25.7 J	ND	ND
7440-23-5	SODIUM	20000	ug/l	167000 J	712000 J	32700000 J
7440-28-0	THALLIUM	0.5 (G)	ug/l	5.2 J	4.02 J	ND
7440-62-2	VANADIUM		ug/l	438 J	77.8 J	ND
7440-66-6	ZINC	2000 (G)	ug/l	5910 J	715 J	46.8 J
57-12-5	CYANIDE	200	ug/L	224	126	174
7420 00 5	DISSOLVED METALS		/1		17.1.1	
7429-90-5	ALUMINUM		ug/L		17.1 J	
7440-36-0	ANTIMONY	3	ug/L		9.22 J	
7440-38-2	ARSENIC	25	ug/L		6 J	
	BARIUM	1000	ug/L		243 ND	
7440-41-7	BERYLLIUM	3 (G)	ug/L		ND	
7440-43-9	CADMIUM	5	ug/L		R	
7440-70-2	CALCIUM		ug/L		143000 J	
7440-47-3	CHROMIUM	50	ug/L		7.84 J	
7440-48-4	COBALT		ug/L		ND	
7440-50-8	COPPER	200	ug/L		55.2 J	
7439-89-6	IRON	300	ug/L		60.5	
	LEAD	25	ug/L		3.29 J	
	MAGNESIUM	35000 (G)	ug/L		21200 J	
7439-96-5	MANGANESE	300	ug/L		488 J	
7439-97-6	MERCURY	0.7	ug/L		ND	
7440-02-0	NICKEL	100	ug/L		10.2 J	
7440-09-7	POTASSIUM		ug/L		8920	
7782-49-2	SELENIUM	10	ug/L		16.2	
7440-23-5	SODIUM	20000	ug/L		753000 J	
7440-62-2	VANADIUM		ug/L		ND	
7440-66-6	ZINC	2000 (G)	ug/L		37.6 J	
Notes:		·				

Indicates concentration exceeds standard or guidance value.

(G)

Indicates guidance value.

No standard or guidance value available. NS Indicates compound was not detected. Indicates an estimated concentration. ND

ug/L Micrograms per liter

Remedial Investigation - October 2012 through April 2013

Consolidated	d Edison		Location ID:	MW-7	MW-7	MW-8
	Ludlow Street Site		Sample ID:	MW-7	MW-7-20130404	MW-8
	roundwater Analytical Data		Lab Sample Id:		E1768-12	D4947-02
	mpound Summary		Source:	CTECH	CTECH	CTECH
		NYSDEC	SDG:	D4947	E1768	D4947
		Class GA	Matrix:	WATER	WATER	WATER
		Groundwater	Sampled:	11/27/2012	4/4/2013	11/26/2012
			Validated:	1/23/2013	4/30/2013	1/23/2013
CAS NO.	COMPOUND	Values (1)	UNITS:			
	VOLATILES					
67-64-1	ACETONE	5	ug/l	ND	ND	ND
74-83-9	BROMOMETHANE	5	ug/l	ND	ND	ND
110-82-7	CYCLOHEXANE		ug/l	2.7 J	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	5	ug/l	24	9.8	ND
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	50 (G)	ug/l	ND	ND	ND
108-10-1	METHYL ISOBUTYL KETONE		ug/l	ND	ND	ND
108-87-2	METHYLCYCLOHEXANE		ug/l	3.5 J	0.44 J	ND
100-42-5	STYRENE	5	ug/l	18	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	10	ug/l	ND	1.6 J	0.52 J
	BTEX	10		- 1.2	1.0 0	
71-43-2	BENZENE	1	ug/l	690	1200	ND
100-41-4	ETHYLBENZENE	5	ug/l	270	90	ND
108-88-3	TOLUENE	5	ug/l	270	62.8	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	5	ug/l	420	180	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	5	ug/l	190	130	ND
70 17 0	CVOCs		ug/1	170	150	1,2
74-87-3	CHLOROMETHANE	5	ug/l	ND	ND	ND
156-59-2	CIS-1,2-DICHLOROETHYLENE	5	ug/l	ND	ND	ND
127-18-4	TETRACHLOROETHYLENE(PCE)	5	ug/l	ND	ND	ND
127 10 1	TETRACTEOROETHTEE(TCE)	3	ug/1	TVD	TVD	T\D
	Total VOCs			1888.2	1674.64	0.52
	SEMIVOLATILES					
98-86-2	ACETOPHENONE		ug/l	5.9 J	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	5	ug/l	7.3 J	25.7	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	5	ug/l	ND	ND	ND
86-74-8	CARBAZOLE		ug/l	ND	5.5 J	ND
132-64-9	DIBENZOFURAN		ug/l	ND	6.3 J	ND
91-94-1	3,3'-DICHLOROBENZIDINE	5	ug/l	ND	ND	ND
99-09-2	3-NITROANILINE	5	ug/l	ND	ND	ND
108-95-2	PHENOL	1	ug/l	ND	ND	ND
	PAHs					
83-32-9	ACENAPHTHENE	20 (G)	ug/l	12.1	47.1	ND
208-96-8	ACENAPHTHYLENE		ug/l	7.9 J	38.6	ND
120-12-7	ANTHRACENE	50 (G)	ug/l	ND	26.2	ND
56-55-3	BENZO(A)ANTHRACENE	0.002 (G)	ug/l	ND	14.8	ND
50-32-8	BENZO(A)PYRENE	ND	ug/l	ND	9.7 J	ND
205-99-2	BENZO(B)FLUORANTHENE	0.002 (G)	ug/l	ND	8.2 J	ND
218-01-9	CHRYSENE	0.002 (G)	ug/l	ND	14.6	ND
206-44-0	FLUORANTHENE	50 (G)	ug/l	ND	18.8	ND
86-73-7	FLUORENE	50 (G)	ug/l	9.9 J	46.3	ND
91-57-6	2-METHYLNAPHTHALENE		ug/l	130	340	ND
91-20-3	NAPHTHALENE	10 (G)	ug/l	1300	1600	ND
85-01-8	PHENANTHRENE	50 (G)	ug/l	14.2	110	ND
129-00-0	PYRENE	50 (G)	ug/l	ND	37	ND
	Total PAHs			1474.1	2311.3	ND
I	1		1			
	Total SVOCs			1487.3	2348.8	ND

Notes:

Indicates concentration exceeds standard or guidance value.

(G) Indicates guidance value.

No standard or guidance value available. Indicates compound was not detected. NS ND Indicates an estimated concentration.

Micrograms per liter ug/L

Remedial Investigation - October 2012 through April 2013

C 1: 1. (1.P.1'		Ir ID. I	1007.7	1007.7	MANA
Consolidated			Location ID:	MW-7 MW-7	MW-7 MW-7-20130404	MW-8 MW-8
Ludlow Street Site Validated Groundwater Analytical Data			Sample ID:		E1768-12	MW-8 D4947-02
	-		Lab Sample Id: Source:	D4947-08		
Detected Col	mpound Summary	NYSDEC	Source: SDG:	CTECH D4047	CTECH E1769	CTECH D4047
		Class GA	Matrix:	D4947	E1768	D4947
				WATER	WATER	WATER
		Groundwater	Sampled:	11/27/2012	4/4/2013	11/26/2012
	T		Validated:	1/23/2013	4/30/2013	1/23/2013
CAS NO.	COMPOUND	Values (1)	UNITS:			
	VOLATILES					
	INORGANICS			260.7	244.7	1060 7
7429-90-5	ALUMINUM		ug/l	268 J	244 J	1860 J
7440-36-0	ANTIMONY	3	ug/l	6.86 J	ND	5.06 J
7440-38-2	ARSENIC	25	ug/l	ND	ND	4.85 J
7440-39-3	BARIUM	1000	ug/l	109 J	192	104 J
	BERYLLIUM	3 (G)	ug/l	ND	ND	ND
7440-43-9	CADMIUM	5	ug/l	ND	R	0.3 J
7440-70-2	CALCIUM		ug/l	105000 J	123000 J	78800 J
7440-47-3	CHROMIUM, TOTAL	50	ug/l	ND	7.73 J	ND
7440-48-4	COBALT		ug/l	ND	ND	4.32 J
7440-50-8	COPPER	200	ug/l	1.44 J	2.06 J	20.7 J
7439-89-6	IRON	300	ug/l	ND	1300	3790 J
7439-92-1	LEAD	25	ug/l	1.72 J	ND	24.5 J
7439-95-4	MAGNESIUM	35000 (G)	ug/l	37800 J	42200 J	25200 J
7439-96-5	MANGANESE	300	ug/l	1030 J	688 J	1520 J
7439-97-6	MERCURY	0.7	ug/l	ND	ND	ND
7440-02-0	NICKEL	100	ug/l	ND	9 J	ND
7440-09-7	POTASSIUM		ug/l	8420 J	26900	8280 J
7782-49-2	SELENIUM	10	ug/l	ND	ND	ND
7440-22-4	SILVER	50	ug/l	ND	ND	ND
7440-23-5	SODIUM	20000	ug/l	535000 J	2230000 J	628000 J
7440-28-0	THALLIUM	0.5 (G)	ug/l	ND	ND	ND
7440-62-2	VANADIUM		ug/l	ND	ND	5.36 J
7440-66-6	ZINC	2000 (G)	ug/l	ND	10.3 J	ND
57-12-5	CYANIDE	200	ug/L	27	145	ND
	DISSOLVED METALS					
7429-90-5	ALUMINUM		ug/L			
7440-36-0	ANTIMONY	3	ug/L			
7440-38-2	ARSENIC	25	ug/L			
	BARIUM	1000	ug/L			
7440-41-7	BERYLLIUM	3 (G)	ug/L			
7440-43-9	CADMIUM	5	ug/L			
7440-70-2	CALCIUM		ug/L			
7440-47-3	CHROMIUM	50	ug/L			
7440-48-4	COBALT		ug/L			
7440-50-8	COPPER	200	ug/L			
7439-89-6	IRON	300	ug/L			
	LEAD	25	ug/L			
	MAGNESIUM	35000 (G)	ug/L			
7439-96-5	MANGANESE	300	ug/L			
7439-97-6	MERCURY	0.7	ug/L			
7440-02-0	NICKEL	100	ug/L			
7440-09-7	POTASSIUM		ug/L			
7782-49-2	SELENIUM	10	ug/L			
7440-23-5	SODIUM	20000	ug/L			
7440-62-2	VANADIUM		ug/L			
7440-66-6	ZINC	2000 (G)	ug/L			
Notes:						

Indicates concentration exceeds standard or guidance value.

(G)

Indicates guidance value.

No standard or guidance value available. NS Indicates compound was not detected. Indicates an estimated concentration. ND

ug/L Micrograms per liter

Table 4 **Summary of Groundwater Analytical Data** Former Ludlow Street Works Site

Consolidated Edison Company of New York Remedial Investigation - October 2012 through April 2013

Ladlow Street Site	Consolidated	1 Edison		Location ID:	MW-18	MW-9	MW-9
Validated Groundwater Analytical Data Lab Sample left D4947-03 D4947-04 E1768-01 CTECH C							
Detected Compound Summary							
NYSDEC Clas GA Groundwater Clas GA Groundwater Standar/Gridiance Martix WATER WATER		•					
Class GA Groundwater Standards/Guidance WATER WATER Standards/Guidance Validance Valid	Detected Co.	inpound Summary	NYSDEC				
Cas No. Compound Cas No. Compound Cas No. Ca							
Standards/Gnidance Validated: 1/23/2013 1/23/2013 4/30/2							
CAS.NO. COMPOUND							
Total Part Color	CAS NO	COMPOLIND			172372013	1/23/2013	4/30/2013
10-84-1 ACETONE	CAS NO.		v aiues	UNIIS.			
74-83-9 BROMOMETHANE 5 ug/l ND ND ND ND	67-64-1		5	119/1	ND	ND	ND
110-82-7							
98-82-8 ISOPROPYLBENZENE (CUMENE) 5 ug/l ND ND ND ND ND ND ND N							
Table Tabl				-			
108-10-1 METHYL ISOBUTYL RETONE		· · · · · · · · · · · · · · · · · · ·		-			
108-87-2 METHYLCYCLOHEXANE ug/l ND ND ND ND ND 1034-04-4 TERT-BUTYL METHYL ETHER 10 ug/l 0.49 J 1.4 J 1.6 J 104-25 BTEX		· · · · · · · · · · · · · · · · · · ·	` ′				
100-42-5 STYRENE 5 ug/l 0.49 J 1.4 J 1.6 J							
1634-04-4 TERT-BUTYL METHYL ETHER				-			
BTEX				-			
1	1034-04-4		10	ug/1	U.47 J	1.+ J	1.0 J
100-41-4 ETHYLBENZENE 5 ug/l ND ND ND ND ND ND ND N	71_43_2		1	119/1	ND	ND	ND
108-88-3 TOLUENE				-			
XYLMP				-			
95-47-6 O-XYLENE (1,2-DIMETHYLBENZENE) 5 ug/l ND ND ND ND				-			
CVOCS				-			
74-87-3 CHLOROMETHANE 5 ug/l ND ND ND ND 156-59-2 CIS-1,2-DICHLOROETHYLENE 5 ug/l ND ND ND ND 127-18-4 TETRACHLOROETHYLENE(PCE) 5 ug/l ND ND ND ND	75-47-0		3	ug/1	ND	ND	ND
156-59-2 CIS-1,2-DICHLOROETHYLENE 5 ug/l ND ND ND ND ND ND ND N	74-87-3		5	110/1	ND	ND	ND
127-18-4 TETRACHLOROETHYLENE(PCE) 5 ug/l ND ND ND ND ND				-			
Total VOCs				-			
SEMIVOLATILES	127 10 1	TETRICIDENCETTT EEL (ECE)	5	ug/1	T\D	ND	TVD
98-86-2 ACETOPHENONE 92-52-4 BIPHENYL (DIPHENYL) 5 ug/l ND ND ND ND 86-74-8 CARBAZOLE ug/l ND ND ND ND 86-74-8 CARBAZOLE ug/l ND ND ND ND 132-64-9 DIBENZOFURAN ug/l ND ND ND ND ND 91-94-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND ND ND 91-94-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND ND ND 108-95-2 PHENOL ug/l ND ND ND ND ND 108-95-2 PHENOL ug/l ND ND ND ND ND 108-95-2 PHENOL ug/l ND ND ND ND ND 108-95-3 ACENAPHTHENE ug/l ND ND ND ND 120-12-7 ANTHRACENE 5 0(G) ug/l ND ND ND ND 50-32-8 BENZO(A)ANTHRACENE ND ug/l ND ND ND ND 50-32-8 BENZO(A)PYRENE ND ug/l ND ND ND 50-32-8 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND ND ND 0.002 (G) ug/l ND ND ND 0.002 (G) ug/l ND ND ND 0.003 (G) ug/l ND ND ND 0.004 (H) ND ND 0.005 (H) ND ND 0.005 (H) ND ND 0.005 (H) ND ND ND 0.005 (H) ND ND ND 0.007 (H) ND ND ND 0.008 (H) ND ND ND 0.009 (H) ND ND ND ND ND 0.009 (H) ND ND ND ND ND 0.009 (H) ND		Total VOCs			0.49	1.4	1.6
92-52-4 BIPHENYL (DIPHENYL) 5 ug/l ND ND ND 117-81-7 BIS(2-ETHYLHEXYL) PHTHALATE 5 ug/l ND ND ND 132-64-8 CARBAZOLE ug/l ND ND ND 132-64-9 DIBENZOFURAN ug/l ND ND ND 91-94-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND 90-90-2 3-NITROANILINE 5 ug/l ND ND ND 108-95-2 PHENOL ug/l ND ND ND 108-95-2 PHENOL ug/l ND ND ND ND 120-12-7 ANTHRACENE 5 Ug/l ND N							
117-81-7 BIS(2-ETHYLHEXYL) PHTHALATE 5 ug/l ND ND ND 86-74-8 CARBAZOLE							
S6-74-8 CARBAZOLE		, , , , , , , , , , , , , , , , , , ,					
132-64-9 DIBENZOFURAN ug/l ND ND ND 91-94-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND 99-09-2 3-NITROANILINE 5 ug/l ND ND ND 108-95-2 PHENOL 1 ug/l ND ND ND PAHS		· · · · · · · · · · · · · · · · · · ·	5				
91-94-1 3,3'-DICHLOROBENZIDINE 5 ug/l ND ND ND ND ND 108-95-2 PHENOL 1 ug/l ND							
99-09-2 3-NITROANILINE 5 ug/l ND ND ND ND 108-95-2 PHENOL 1 ug/l ND				-			
1				-			
PAHS S3-32-9 ACENAPHTHENE 20 (G) ug/l ND ND ND ND ND 208-96-8 ACENAPHTHYLENE ug/l ND ND ND ND ND ND ND N				ug/l			
S3-32-9 ACENAPHTHENE 20 (G) ug/l ND ND ND	108-95-2		1	ug/l	ND	ND	ND
208-96-8 ACENAPHTHYLENE ug/l ND ND ND 120-12-7 ANTHRACENE 50 (G) ug/l ND ND ND 56-55-3 BENZO(A)ANTHRACENE 0.002 (G) ug/l ND ND ND 50-32-8 BENZO(A)PYRENE ND ug/l ND ND ND 205-99-2 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND ND ND 218-01-9 CHRYSENE 0.002 (G) ug/l ND ND ND 206-44-0 FLUORANTHENE 50 (G) ug/l ND ND ND 86-73-7 FLUORENE 50 (G) ug/l ND ND ND 86-73-6 2-METHYLNAPHTHALENE ug/l ND ND ND 91-20-3 NAPHTHALENE 10 (G) ug/l ND ND ND 85-01-8 PHENANTHRENE 50 (G) ug/l ND ND ND 129-00-0 PYRENE 50 (G) ug/l ND ND ND							
120-12-7 ANTHRACENE 50 (G) ug/l ND ND ND ND			20 (G)	-			
56-55-3 BENZO(A)ANTHRACENE 0.002 (G) ug/l ND ND ND 50-32-8 BENZO(A)PYRENE ND ug/l ND ND ND 205-99-2 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND ND ND 218-01-9 CHRYSENE 0.002 (G) ug/l ND ND ND 206-44-0 FLUORANTHENE 50 (G) ug/l ND ND ND 86-73-7 FLUORENE 50 (G) ug/l ND ND ND 91-57-6 2-METHYLNAPHTHALENE ug/l ND ND ND 91-20-3 NAPHTHALENE 10 (G) ug/l ND ND ND 85-01-8 PHENANTHRENE 50 (G) ug/l ND ND ND 129-00-0 PYRENE 50 (G) ug/l ND ND ND ND ND ND ND ND ND				-			
50-32-8 BENZO(A)PYRENE ND ug/l ND ND ND 205-99-2 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND ND ND 218-01-9 CHRYSENE 0.002 (G) ug/l ND ND ND 206-44-0 FLUORANTHENE 50 (G) ug/l ND ND ND 86-73-7 FLUORENE 50 (G) ug/l ND ND ND 91-57-6 2-METHYLNAPHTHALENE ug/l ND ND ND 91-20-3 NAPHTHALENE 10 (G) ug/l ND ND ND 85-01-8 PHENANTHRENE 50 (G) ug/l ND ND ND 129-00-0 PYRENE 50 (G) ug/l ND ND ND ND ND ND ND ND				-			
205-99-2 BENZO(B)FLUORANTHENE 0.002 (G) ug/l ND ND ND ND		· ·	` /	-			
218-01-9 CHRYSENE 0.002 (G) ug/l ND ND ND 206-44-0 FLUORANTHENE 50 (G) ug/l ND ND ND 86-73-7 FLUORENE 50 (G) ug/l ND ND ND 91-57-6 2-METHYLNAPHTHALENE ug/l ND ND ND 91-20-3 NAPHTHALENE 10 (G) ug/l ND ND ND 85-01-8 PHENANTHRENE 50 (G) ug/l ND ND ND 129-00-0 PYRENE 50 (G) ug/l ND ND ND Total PAHs ND ND ND ND ND ND ND				-			
206-44-0 FLUORANTHENE 50 (G) ug/l ND ND ND ND ND ND ND N		` /		-			
86-73-7 FLUORENE 50 (G) ug/l ND ND ND 91-57-6 2-METHYLNAPHTHALENE ug/l ND ND ND 91-20-3 NAPHTHALENE 10 (G) ug/l ND ND ND 85-01-8 PHENANTHRENE 50 (G) ug/l ND ND ND 129-00-0 PYRENE 50 (G) ug/l ND ND ND Total PAHs			* *	-			
91-57-6 2-METHYLNAPHTHALENE ug/l ND ND ND ND 91-20-3 NAPHTHALENE 10 (G) ug/l ND				-			
91-20-3 NAPHTHALENE 10 (G) ug/l ND			* *	-			
85-01-8				-			
129-00-0 PYRENE 50 (G) ug/l ND ND ND ND ND ND ND N	91-20-3			-			
Total PAHs ND ND ND				-			
	129-00-0	PYRENE	50 (G)	ug/l	ND	ND	ND
Total SVOCs ND ND ND		Total PAHs			ND	ND	ND
		Total SVOCs			ND	ND	ND

Notes:

Indicates concentration exceeds standard or guidance value.

(G) Indicates guidance value.

No standard or guidance value available. Indicates compound was not detected. NS ND Indicates an estimated concentration.

Micrograms per liter ug/L

Remedial Investigation - October 2012 through April 2013

Consolidated	Edison		Location ID:	MW-18	MW-9	MW-9
Ludlow Street Site			Sample ID:	MW-18	MW-9	MW-9-20130405
Validated Groundwater Analytical Data			Lab Sample Id:		D4947-04	E1768-01
	mpound Summary		Source:	CTECH	CTECH	CTECH
Detected Co.	inpound Summary	NYSDEC	SDG:	D4947	D4947	E1768
		Class GA	Matrix:	WATER	WATER	WATER
		Groundwater	Sampled:	11/26/2012	11/26/2012	4/5/2013
			Validated:	1/23/2013	1/23/2013	4/30/2013
CAS NO.	COMPOUND	Values (1)	UNITS:	172372013	1/23/2013	1/30/2013
CAS NO.	VOLATILES	values	UNITS.			
	INORGANICS					
7429-90-5	ALUMINUM		ug/l	9290 J	ND	91.9 J
7440-36-0	ANTIMONY	3	ug/l	ND	ND	ND
7440-38-2	ARSENIC	25	ug/l	15.3 J	13.7	9.05 J
7440-39-3	BARIUM	1000	ug/l	169 J	478 J	380
7440-41-7	BERYLLIUM	3 (G)	ug/l	0.38 J	ND	ND
7440-43-9	CADMIUM	5	ug/l	0.87 J	0.27 J	R
7440-70-2	CALCIUM		ug/l	90000 J	198000 J	196000 J
7440-47-3	CHROMIUM, TOTAL	50	ug/l	ND	ND	ND
7440-48-4	COBALT		ug/l	12.8 J	ND	ND ND
7440-50-8	COPPER	200	ug/l	59 J	1.11 J	ND
7439-89-6	IRON	300	ug/l	18700 J	17400 J	11300
7439-92-1	LEAD	25	ug/l	85.2 J	ND	ND
7439-95-4	MAGNESIUM	35000 (G)	ug/l	31000 J	88100 J	77600 J
7439-96-5	MANGANESE	300	ug/l	2030 J	4210 J	3240 J
7439-97-6	MERCURY	0.7	ug/l	ND	ND	ND
7440-02-0	NICKEL	100	ug/l	ND	ND	ND
7440-02-0	POTASSIUM		ug/l	10900 J	22200 J	15100
7782-49-2	SELENIUM	10	ug/l	ND	2.82 J	ND
7440-22-4	SILVER	50	ug/l	ND	ND	ND
7440-23-5	SODIUM	20000	ug/l	677000 J	ND	1210000 J
7440-28-0	THALLIUM	0.5 (G)	ug/l	ND	ND	3.46 J
7440-62-2	VANADIUM		ug/l	20.6 J	ND	ND
7440-66-6	ZINC	2000 (G)	ug/l	90.7 J	ND	ND
57-12-5	CYANIDE	200	ug/L	ND	ND	3 J
0, 120	DISSOLVED METALS	200	48,2	1,2	112	2 0
7429-90-5	ALUMINUM		ug/L			
7440-36-0	ANTIMONY	3	ug/L			
7440-38-2	ARSENIC	25	ug/L			
7440-39-3	BARIUM	1000	ug/L			
7440-41-7	BERYLLIUM	3 (G)	ug/L			
7440-43-9	CADMIUM	5	ug/L			
7440-70-2	CALCIUM		ug/L			
7440-47-3	CHROMIUM	50	ug/L			
7440-48-4	COBALT		ug/L			
7440-50-8	COPPER	200	ug/L			
7439-89-6	IRON	300	ug/L			
7439-92-1	LEAD	25	ug/L			
7439-95-4	MAGNESIUM	35000 (G)	ug/L			
7439-96-5	MANGANESE	300	ug/L			
7439-97-6	MERCURY	0.7	ug/L			
7440-02-0	NICKEL	100	ug/L			
7440-09-7	POTASSIUM		ug/L			
7782-49-2	SELENIUM	10	ug/L			
7440-23-5	SODIUM	20000	ug/L			
7440-62-2	VANADIUM		ug/L			
7440-66-6	ZINC	2000 (G)	ug/L			
Notes:		=====(0)				1

Notes:

Indicates concentration exceeds standard or guidance value.

(G)

Indicates guidance value.

No standard or guidance value available. NS ND Indicates compound was not detected. Indicates an estimated concentration.

ug/L Micrograms per liter

Table 5 **Summary of Soil Gas Analytical Results Former Ludlow Street Works** Consolidtaed Edison Company of New York

Remedial Investigation - October 2012 through April 2013

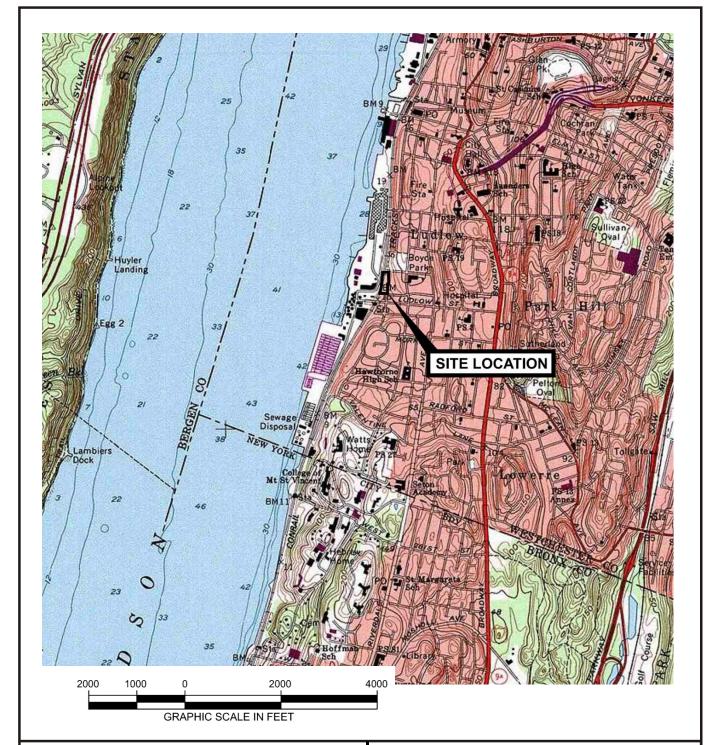
Consolidated Edisor	n	Location ID:	Target	SV-1	SV-3	SV-4
Former Ludlow Stre		Sample ID:	Soil Gas	SV-1 (7')	SV-3 (8')	SV-4 (8')
Validated Air Analy	rtical Data	Lab Sample Id:	Concentration	1210554-03A	1210554-01A	1210554-02A
Detected Compound	d Summary	Source:	(1)(2)	ATL	ATL	ATL
•	•	SDG:		1210554	1210554	1210554
		Matrix:		Soil Gas	Soil Gas	Soil Gas
		Sampled:		10/24/2012	10/24/2012	10/24/2012
		Validated:		4/15/2013	4/15/2013	4/15/2013
CAS NO.	COMPOUND	UNITS:				
	CHLORINATED VOC					
71-55-6	1,1,1-Trichloroethane	UG/M3	22,000	6.5	ND	ND
56-23-5	Carbon tetrachloride	UG/M3	16	ND	ND	ND
156-59-2	cis-1,2-Dichloroethene	UG/M3	350	ND	ND	ND
127-18-4	Tetrachloroethene	UG/M3	81	28	100	24
79-01-6	Trichloroethene	UG/M3	2.2	ND	ND	ND
	NON-MGP RELATED VOC					
106-99-0	1,3-Butadiene	UG/M3	NS	34	ND	4.4
541-73-1	1,3-Dichlorobenzene	UG/M3	1,100	ND	ND	ND
106-46-7	1,4-Dichlorobenzene	UG/M3	8,000	ND	ND	ND
540-84-1	2,2,4-Trimethylpentane	UG/M3	NS	12	ND	ND
78-93-3	2-Butanone (Methyl Ethyl Ketone)	UG/M3	10,000	11	ND	ND
591-78-6	2-Hexanone	UG/M3	NS	ND	ND	ND
107-83-5	2-Methylpentane	UG/M3	NS	ND	ND	ND
67-63-0	2-Propanol	UG/M3	NS	ND	ND	ND
108-10-1	4-Methyl-2-pentanone	UG/M3	800	ND	ND	ND
67-64-1	Acetone	UG/M3	3,500	260	150	220
75-27-4	Bromodichloromethane	UG/M3	14	ND	ND	ND
74-83-9	Bromomethane	UG/M3	50	ND	ND	ND
75-15-0	Carbon disulfide	UG/M3	7,000	17	ND	14
67-66-3	Chloroform	UG/M3	11	ND	ND	ND
74-87-3	Chloromethane	UG/M3	240	ND	ND	ND
98-82-8	Cumene	UG/M3	4,000	ND	ND	ND
110-82-7	Cyclohexane	UG/M3	NS	7.9	ND	ND
64-17-5	Ethanol	UG/M3	NS	ND	ND	ND
75-69-4	Freon 11 (Trichlorofluoromethane)	UG/M3	7,000	ND	ND	ND
76-13-1	Freon 113	UG/M3	300,000	ND	ND	ND
75-71-8	Freon 12 (Dichlorodifluoromethane)	UG/M3	2,000	5.9	ND	ND
142-82-5	Heptane	UG/M3	NS	13	ND	190
110-54-3	Hexane	UG/M3	2,000	27	ND	360
75-09-2	Methylene chloride	UG/M3	520	ND	ND	ND
103-65-1	Propylbenzene	UG/M3	1,400	ND	ND	ND
100-42-5	Styrene	UG/M3	10,000	ND	ND	ND
109-99-9	Tetrahydrofuran	UG/M3	NS	ND	ND	ND
	POSSIBLY MGP RELATED VOC					
95-63-6	1,2,4-Trimethylbenzene	UG/M3	60	ND	ND	ND
108-67-8	1,3,5-Trimethylbenzene	UG/M3	60	ND	ND	ND
622-96-8	4-Ethyltoluene	UG/M3	NS	ND	ND	ND
100-41-4	Ethyl Benzene	UG/M3	220	5.4	ND	5.8
71-43-2	Benzene	UG/M3	31	20	ND	8.6
496-11-7	Indan	UG/M3	NS	ND	ND	ND
78-78-4	Isopentane	UG/M3	NS	ND	ND	ND
108-38-3/106-42-3	m,p-Xylene	UG/M3	70,000	19	11	20
91-20-3	Naphthalene	UG/M3	30	ND	ND	ND
95-47-6	o-Xylene	UG/M3	70,000	5.2	ND	7
108-88-3	Toluene	UG/M3	4,000	35	23	60
			,			

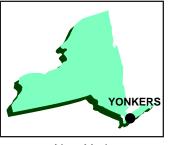
Notes:

^{1.} United States Environmental Protection Agency, "OSWER Draft Guidance for Evaluating the Vapor Intrusion to Indoor Air Pathway from Groundwater and Soils (Subsurface Vapor Intrusion Guidance)", Table 2b, EPA530-D-02-004, November 2002.

^{2.} NS indicates Not Specified.

^{3.} Bold/highlighted results indicate exceedance of target concentration.





New York

N

SOURCE:

MAP CREATED WITH TOPO!TM ©2000 WILDFLOWER PRODUCTIONS (www.topo.com)

FIGURE 1

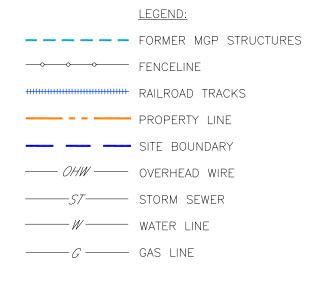
Consolidated Edison Former Ludlow Street Works Yonkers, New York

SITE LOCATION MAP

PARSONS

200 COTTONTAIL LANE, SOMERSET, NJ 08873 PHONE: (732) 537-3500





NOTES:

100

1"=50

SCALE:

- BASE SURVEY DRAWING PROVIDED BY THE CHAZEN COMPANIES. A SURVEY PERFORMED ON THE SITE IN JULY 2010 AND UPDATED IN DECEMBER 2012.
- 2. HISTORIC FEATURES IN BASE DRAWING BASED ON SANBORN FIRE INSURANCE MAPS DATED 1888, 1898, 1917, 1951, & 1991. WESTCHESTER COUNTY ATLAS, 1931., YONKERS TAX ASSESSORS MAP, REVISED 1947 AND FIGURE 3 FROM GEI CONSULTANTS, INC.

FIGURE 2

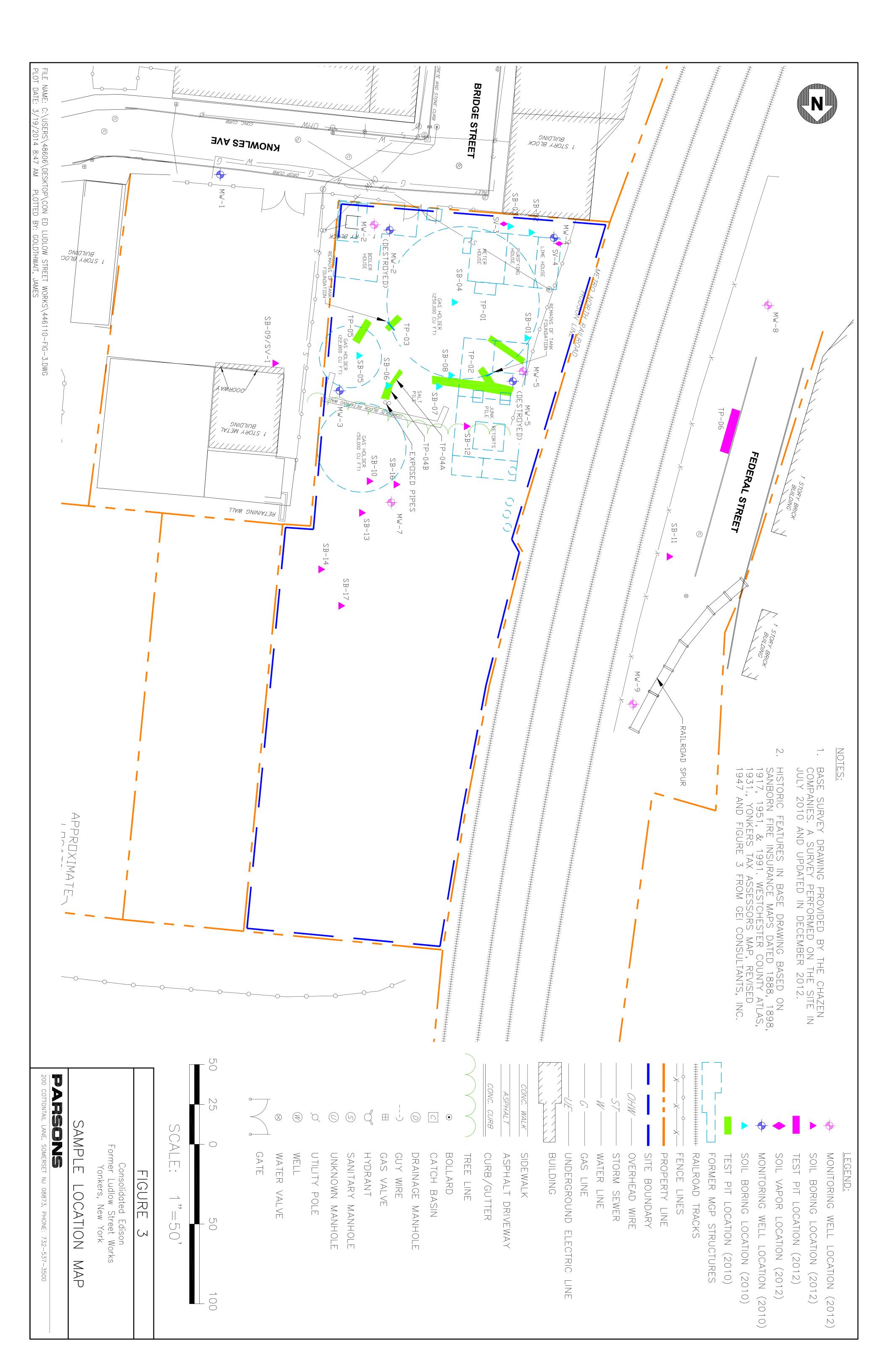
Consolidated Edison Former Ludlow Street Works Yonkers, New York

FORMER MGP STRUCTURES

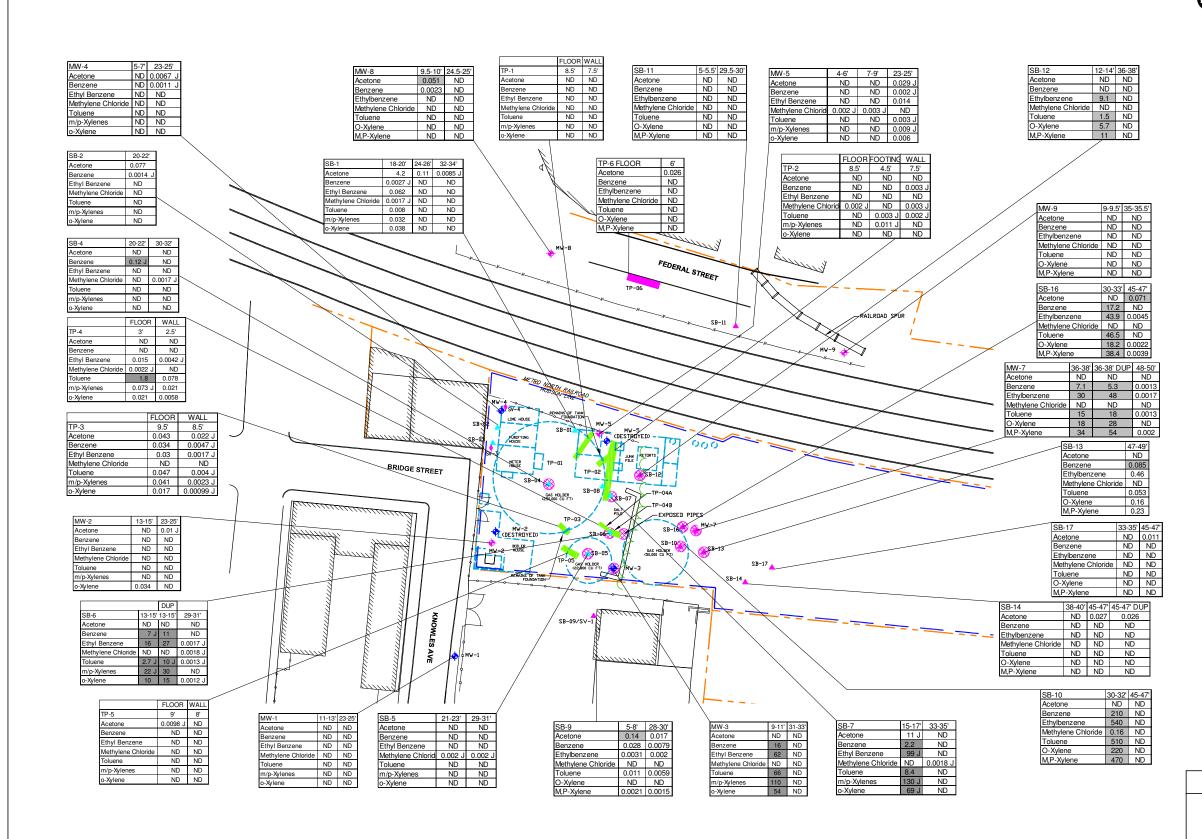
PARSONS

200 COTTONTAIL LANE, SOMERSET NJ 08873, PHONE: 732-537-3500

FILE NAME: C:\USERS\48606\DESKTOP\CON ED LUDLOW STREET WORKS\446110-FIG-2.DWG PLOT DATE: 2/24/2014 10:04 AM PLOTTED BY: GOLDTHWAIT, JAMES







LEGEND:

- ♦ MONITORING WELL LOCATION (2012)
- SOIL BORING LOCATION (2012)
- TEST PIT LOCATION (2012)
- SOIL VAPOR LOCATION (2012)
- MONITORING WELL LOCATION (2010)
- ▲ SOIL BORING LOCATION (2010)
- TEST PIT LOCATION (2010)
- ---- FORMER MGP STRUCTURES
 ---- FENCELINE
- RAILROAD TRACKS
- --- PROPERTY LINE
- - SITE BOUNDARY
- 22222222 BNITDING
 - NAPL OBSERVED

NOTES:

- 1. BASE SURVEY DRAWING PROVIDED BY THE CHAZEN COMPANIES, FROM A SURVEY PERFORMED ON THE SITE IN JULY. 2010 AND UPDATED IN DECEMBER 2012.
- 2. HISTORIC FEATURES IN BASE DRAWING BASED ON SANBORN FIRE INSURANCE MAPS DATED 1888, 1898, 1917, 1951, & 1991. WESTCHESTER COUNTY ATLAS, 1931., YONKERS TAX ASSESSORS MAP, REVISED 1947 AND FIGURE 3 FROM GEI CONSULTANTS, INC.
- 3. ALL CONCENTRATIONS IN MILLIGRAMS PER KILOGRAMS (MG/KG)
- 4. SHADED VALUES EXCEED 6 NYCRR PART 375 UNRESTRICTED SOIL CLEANUP OBJECTIVES (USCO).
- 5. ND NOT DETECTED
- 6. COMPOUNDS DETECTED DURING THE SITE CHARACTERIZATION OR REMEDIAL INVESTIGATION ABOVE USCOs, AT ANY LOCATION, ARE SHOWN.



FIGURE 4

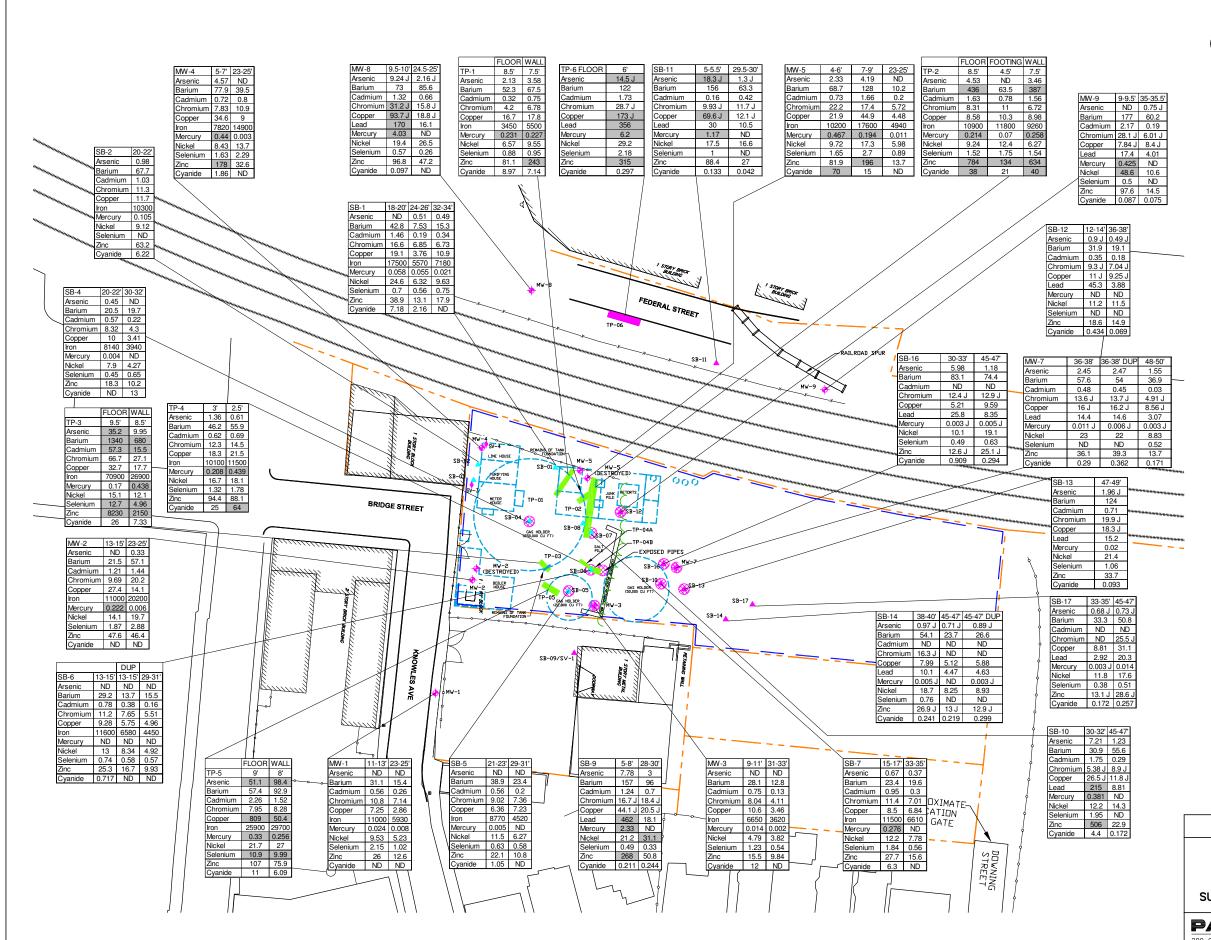
Consolidated Edison Former Ludlow Street Works Yonkers, New York

SUMMARY OF VOC EXCEEDANCES IN SOIL

PARSONS

200 COTTONTAIL LANE, SOMERSET NJ 08873, PHONE: 732-537-3500







LEGEND:

- MONITORING WELL LOCATION (2012)
- SOIL BORING LOCATION (2012)
- TEST PIT LOCATION (2012)
- SOIL VAPOR LOCATION (2012)
- ♦ MONITORING WELL LOCATION (2010)
- SOIL BORING LOCATION (2010)
- TEST PIT LOCATION (2010)
- ---- FORMER MGP STRUCTURES
- RAILROAD TRACKS
- 10 112 110 110 110 10
- ------ PROPERTY LINE
- SITE BOUNDARY
- MILLDING BUILDING

→ FENCELINE

NAPL OBSERVED

NOTES

- . BASE SURVEY DRAWING PROVIDED BY THE CHAZEN COMPANIES, FROM A SURVEY PERFORMED ON THE SITE IN JULY, 2010 AND UPDATED IN DECEMBER 2012.
- HISTORIC FEATURES IN BASE DRAWING BASED ON SANBORN FIRE INSURANCE MAPS DATED 1888, 1898, 1917, 1951, & 1991. WESTCHESTER COUNTY ATLAS, 1931., YONKERS TAX ASSESSORS MAP, REVISED 1947 AND FIGURE 3 FROM GEI CONSULTANTS, INC.
- ALL CONCENTRATIONS IN MILLIGRAMS PER KILOGRAMS (MG/KG)
- 4. SHADED VALUES EXCEED 6 NYCRR PART 375 UNRESTRICTED SOIL CLEANUP OBJECTIVES (USCO).
- 5. ND NOT DETECTED
- 6. COMPOUNDS DETECTED DURING THE SITE CHARACTERIZATION OR REMEDIAL INVESTIGATION ABOVE USCOs, AT ANY LOCATION, ARE SHOWN.



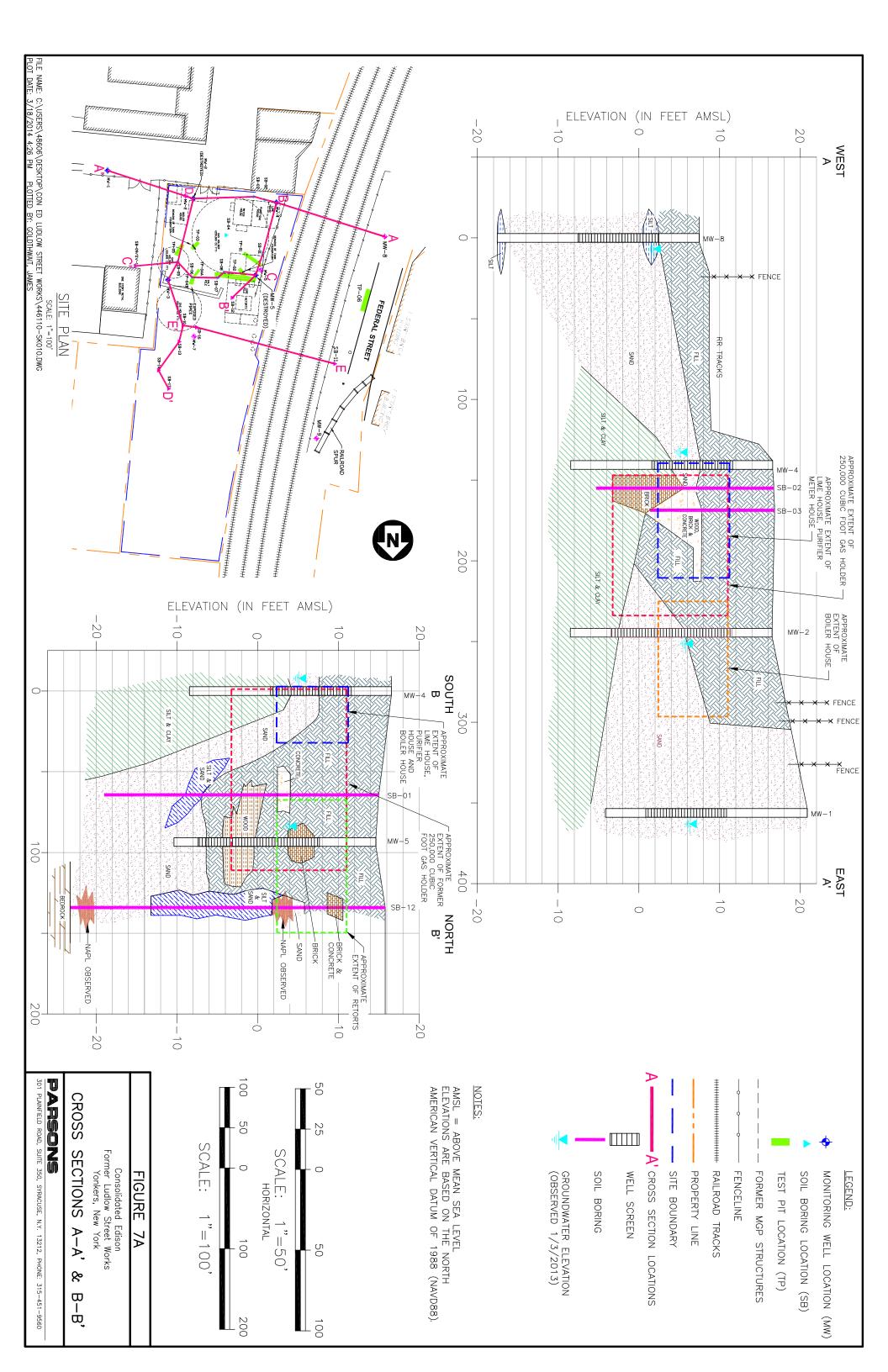
FIGURE 6

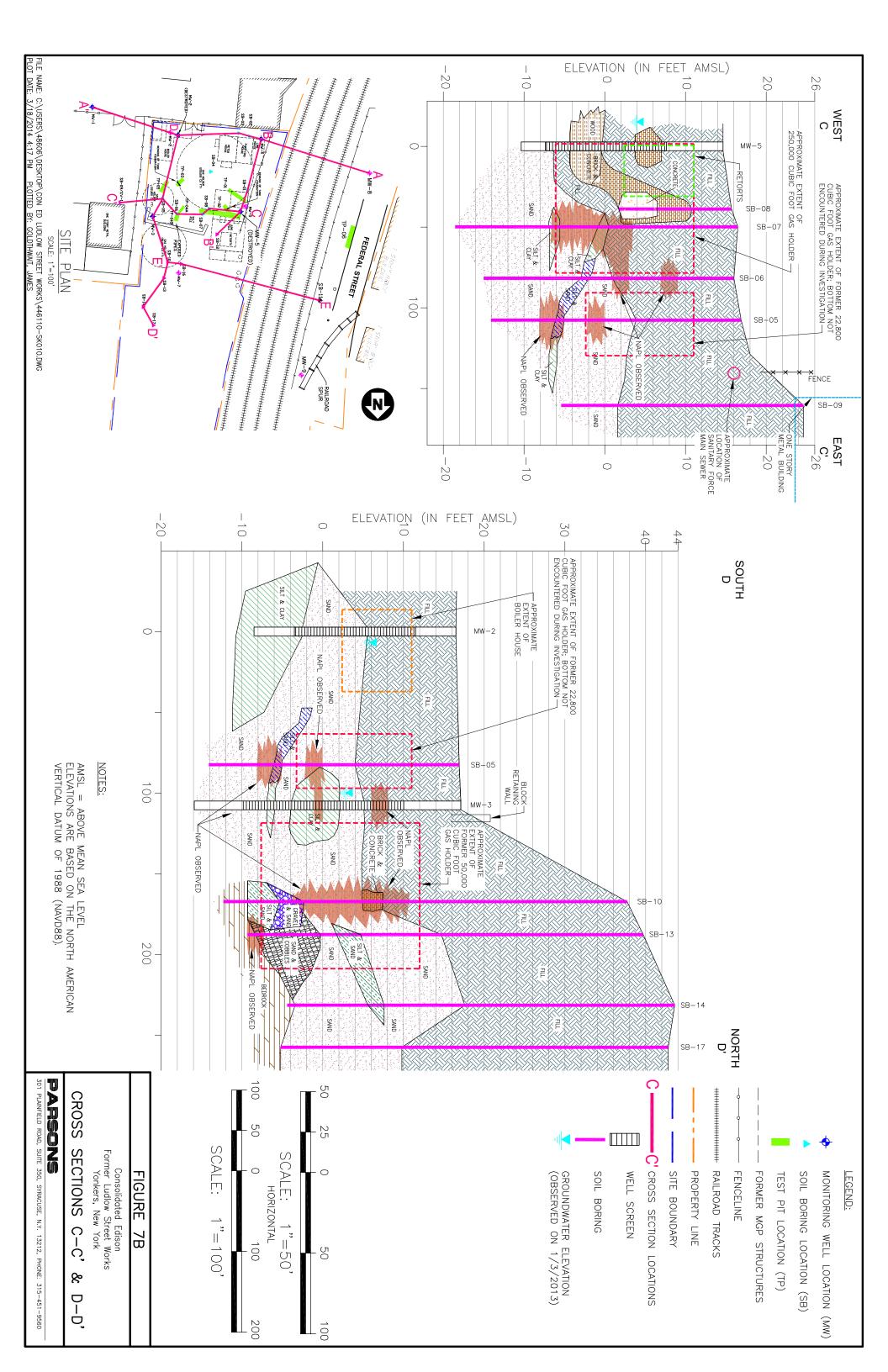
Consolidated Edison Former Ludlow Street Works Yonkers, New York

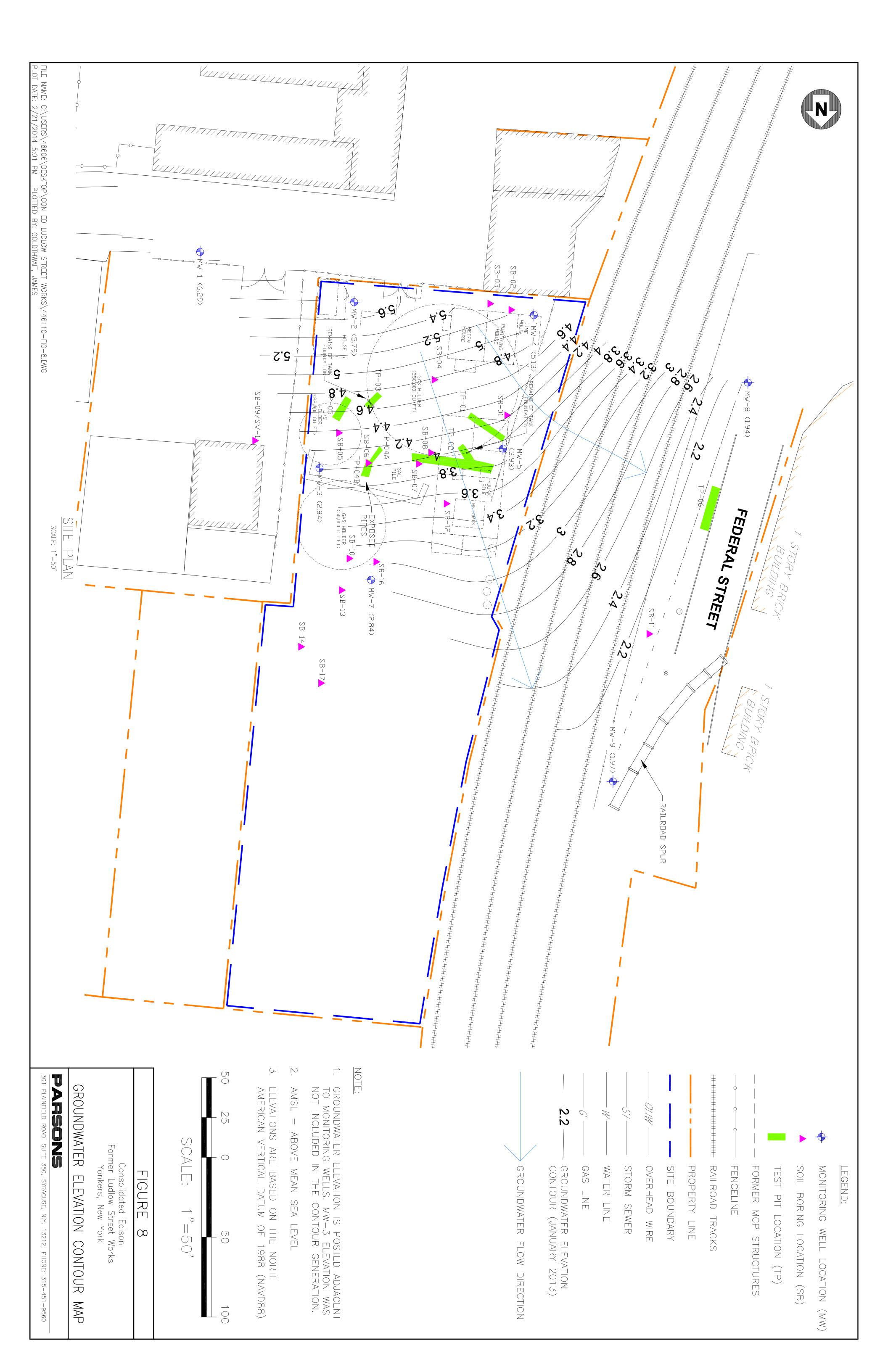
SUMMARY OF INORGANIC EXCEEDANCES IN SOIL

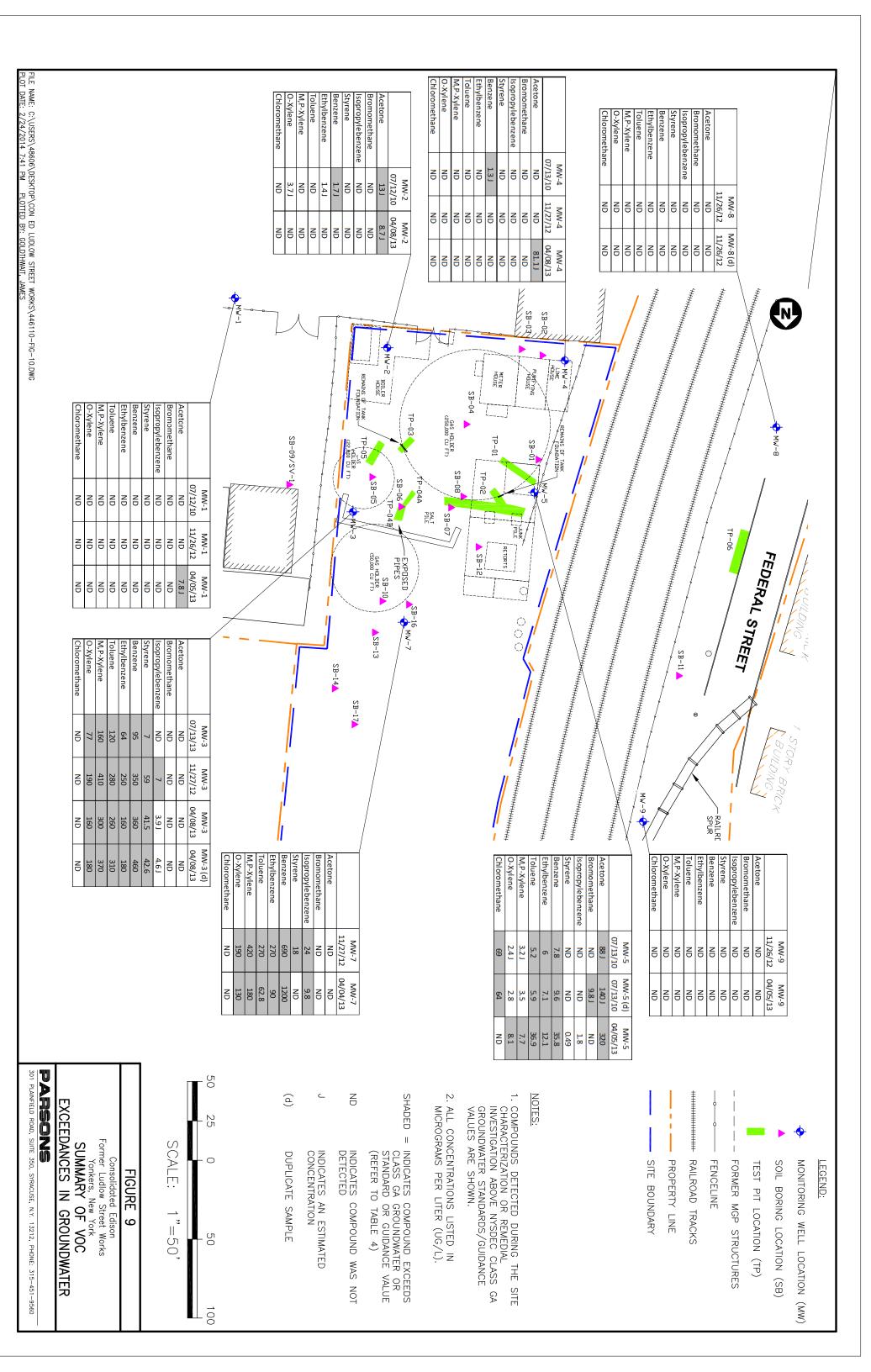
PARSONS

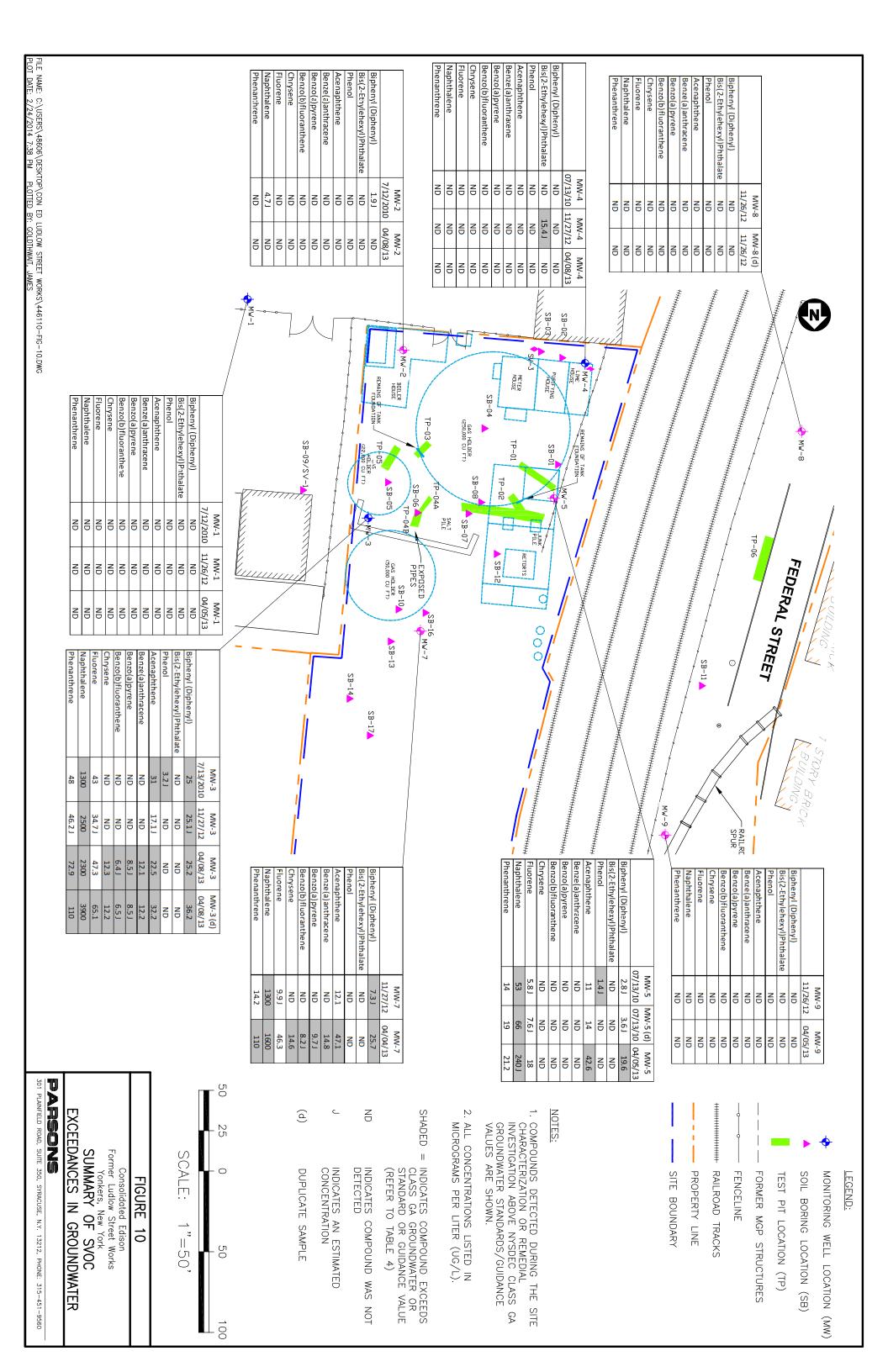
200 COTTONTAIL LANE, SOMERSET NJ 08873, PHONE: 732-537-3500

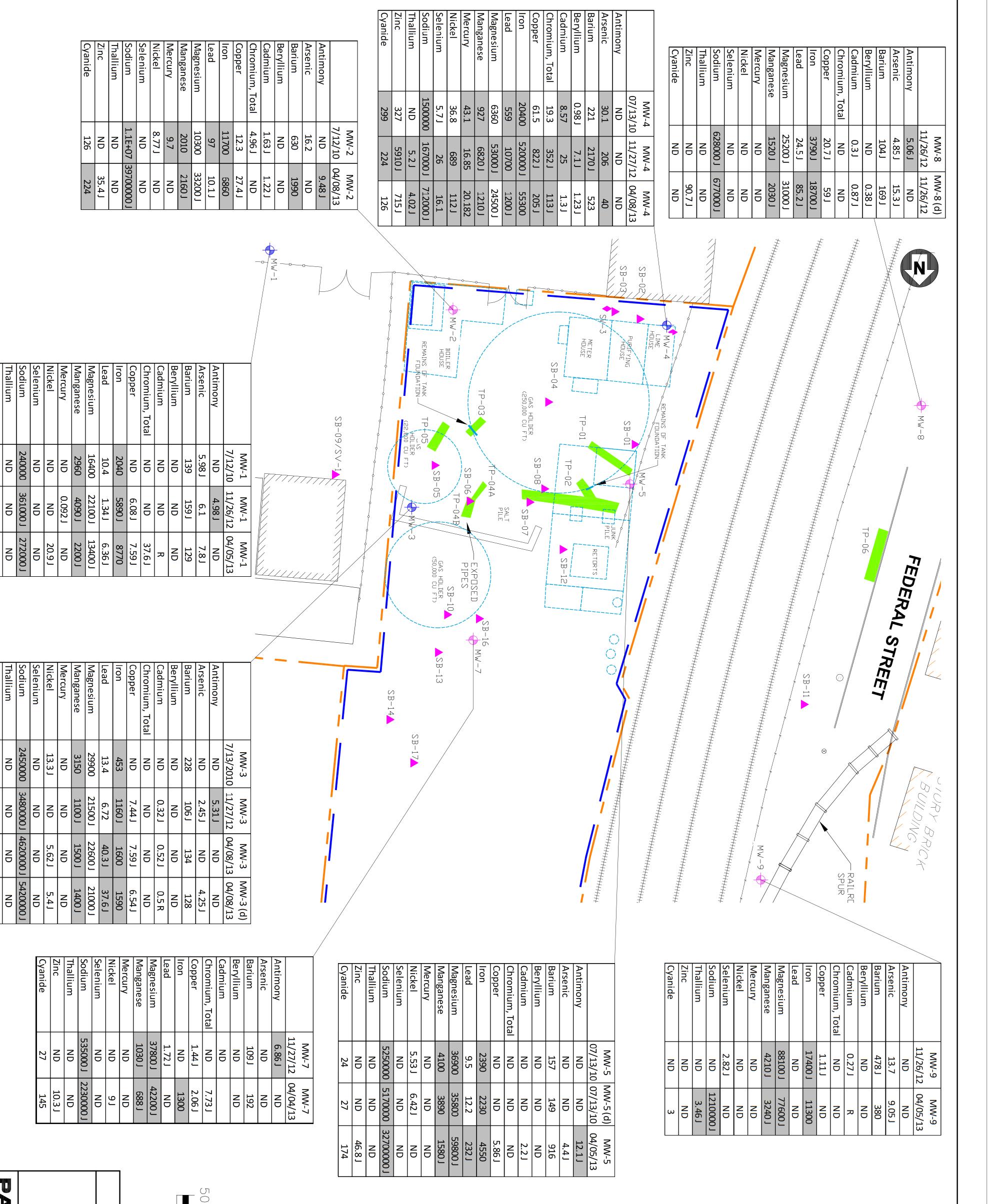












NOTES:

SHE

BOUNDARY

RAILROAD

TRACKS

PROPERTY

FENCELINE

FORMER MGP

STRUCTURES

HES H

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LOCATION (TP)

SOL

BORING

LOCATION (SB)

LEGEND:

MONITORING

 $\forall \exists \Box$

LOCATION

(MM)

COMPOUNDS DETECTED DURING THE SITE CHARACTERIZATION OR REMEDIAL INVESTIGATION ABOVE NYSDEC CLASS GROUNDWATER STANDARDS/GUIDANCE VALUES ARE SHOWN.

 \mathcal{G}

 \mathbb{Z}

INDICATES DETECTED

COMPOUND

WAS

SHADED

INDICATES
CLASS GA
STANDARD
(REFER TO

COMPOUND EXC GROUNDWATER OR GUIDANCE \
D TABLE 4)

VALUE

EXCEEDS

0 7

 $\overline{\bigcirc}$

<u>N</u>

ALL CONCENTRATIONS LISTED IN MICROGRAMS PER LITER (UG/L).

 \subseteq

INDICATES AN ESTIMATED CONCENTRATION

<u>d</u>

DUPLICATE

SAMPLE

25

50

100

S (ALE:

 $\bigcup_{i=1}^{n}$

PLAINFIELD

XCEEDANCES

 \equiv

GROUNDWATER

SUMMARY

INORGANICS

Former Ludlow Yonkers, NMARY OF

 $\mathbb{Z}_{\mathbb{Q}}$

Street W New York

Works

Consolidated

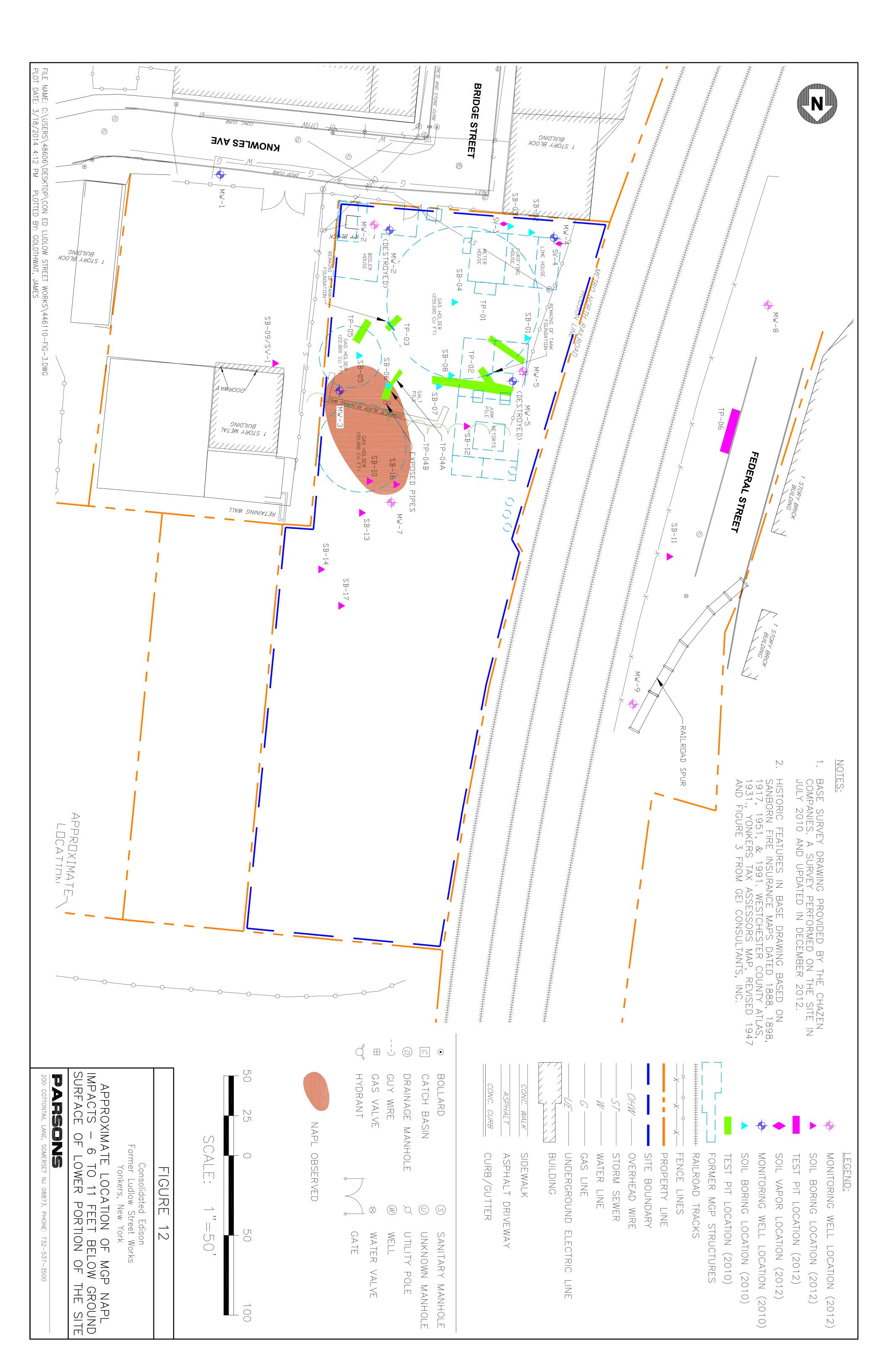
Edison

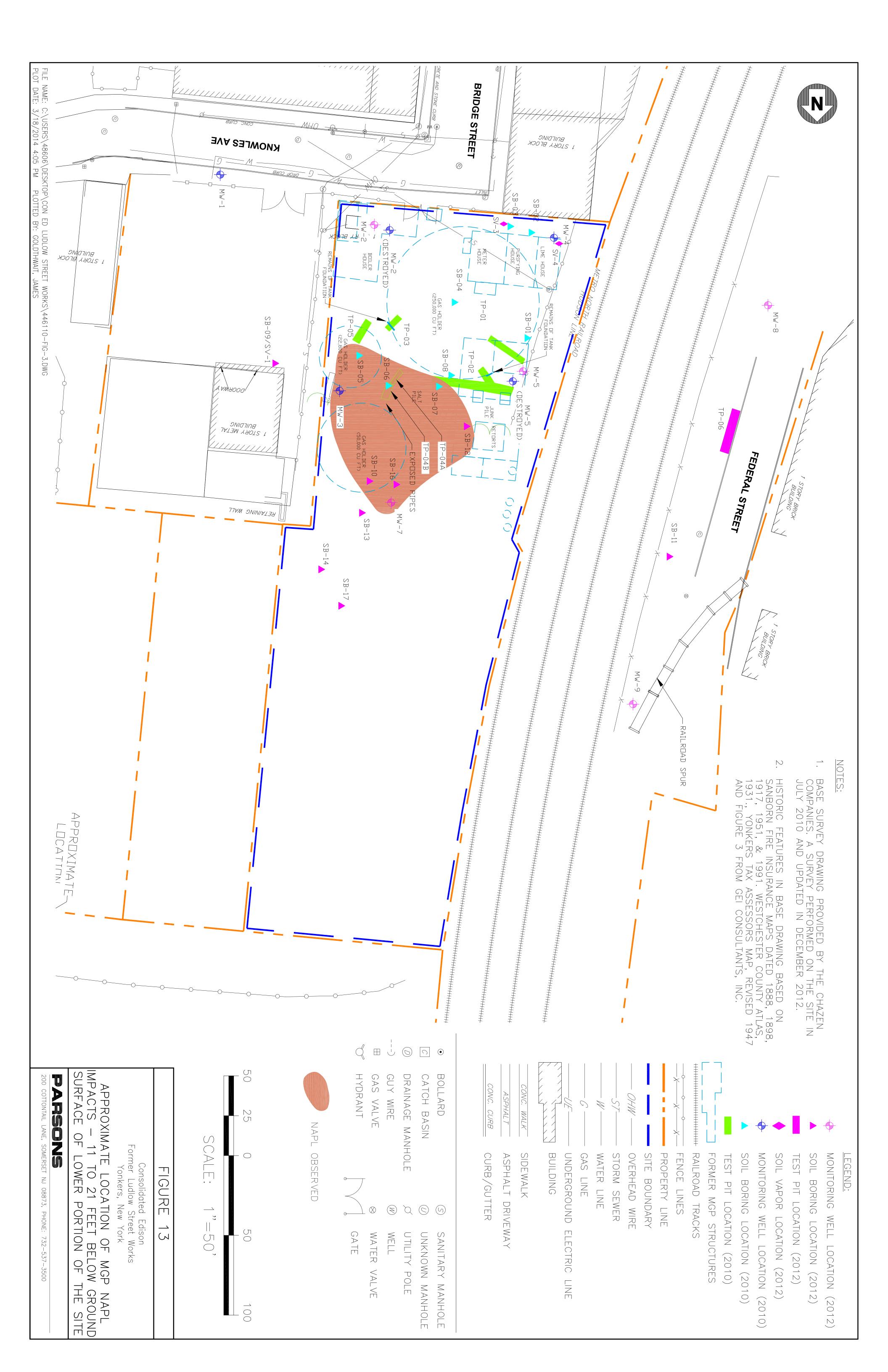
FIGURE

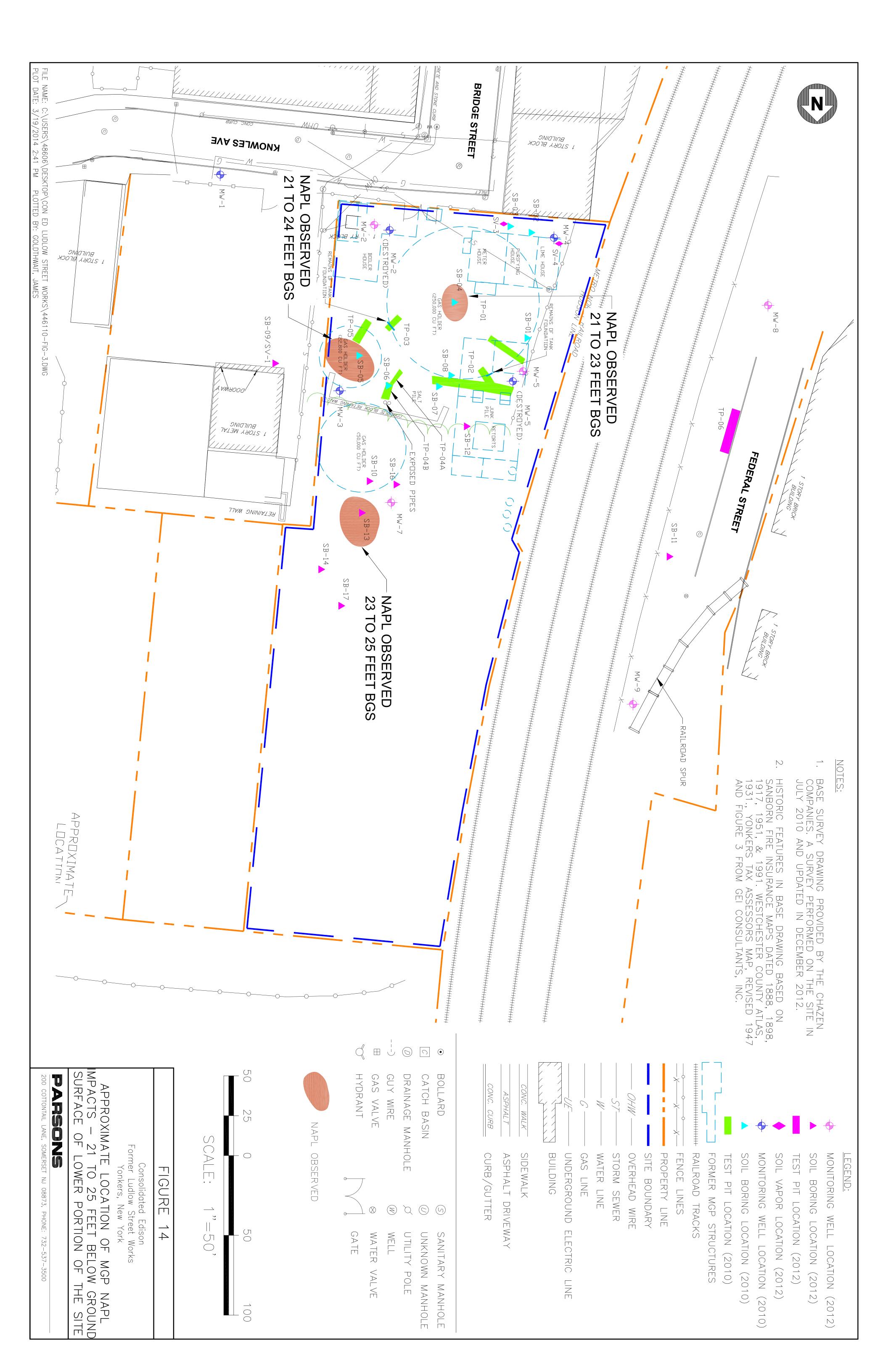
FILE NAME: C:\USERS\48606\DESKTOP\CON ED LUDLOW STREET WORKS\446110-FIG-10.DWGPLOT DATE: 3/19/2014 1:34 PM PLOTTED BY: GOLDTHWAIT, JAMES

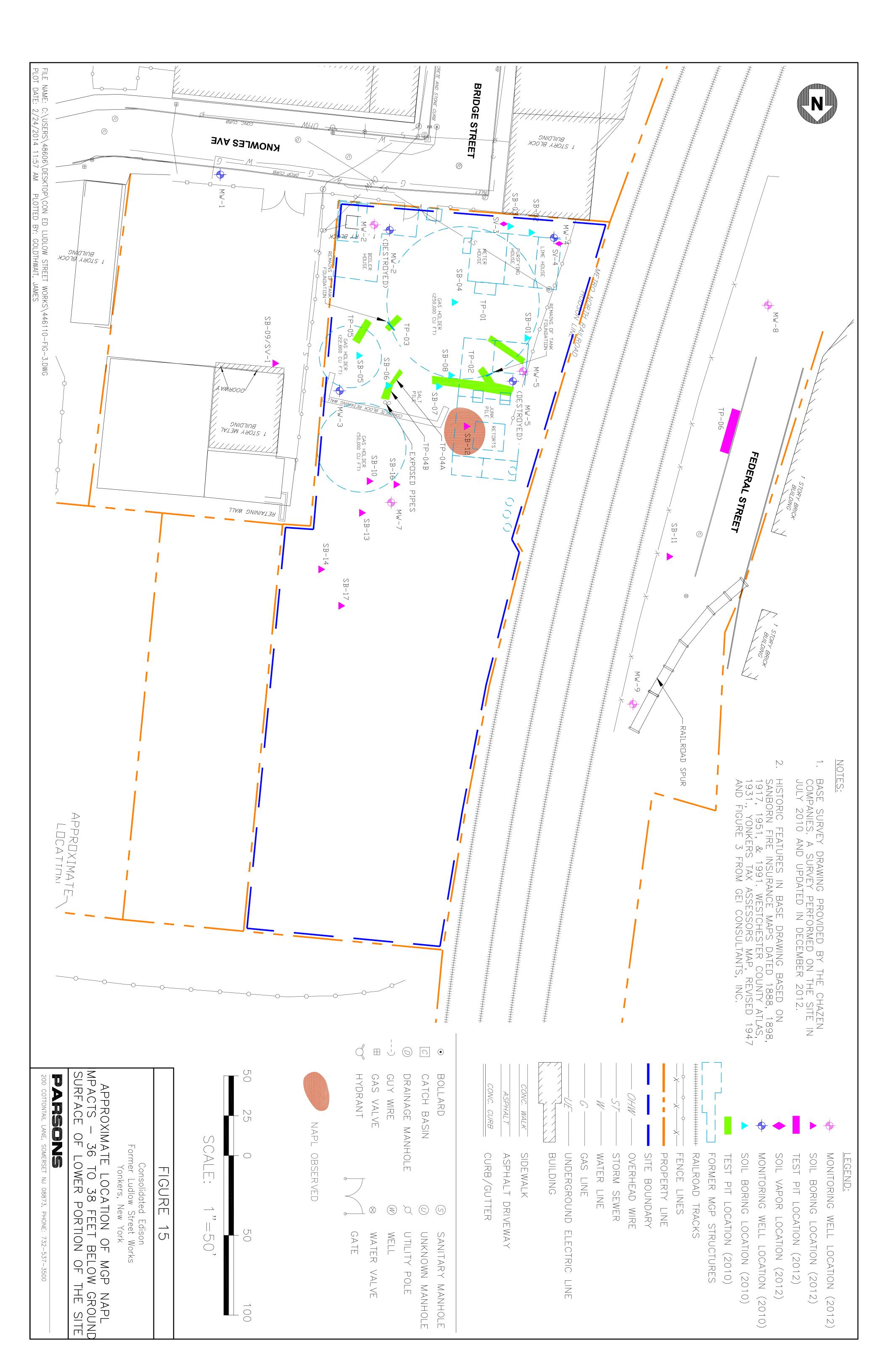
ND

ROAD, 350, SYRACUSE, N.Y. 13212, PHONE: 315-451-9560









	PARSONS TEST PIT RECORD	
PROJECT NAME:	Con Edison: Ludlow Street Works	TEST PIT ID: TP-6
PROJECT NUMBER:	446110	LOCATION:
WEATHER	Partly cloudy, 60s, 5-10 mph E	Approximate L X W X D
DATE/TIME START:	November 7, 2012/0910	= 29' X 4' X 6'
DATE/TIME FINISH:	November 7, 2012/1120	
CONTRACTOR:	ADT	West of MTA Tracks, along Federal Street
INSPECTOR:	Zohar Lavy	
DEPTH (feet bgs)	FIELD IDENTIFICATION OF MATERIAL	COMMENTS
0	0-5' Moist, dark brown/grey fine to coarse SAND and fine to coarse sub-angular to sub-round Gravel, some Slag, little Metal, trace Glass, trace Brick	Naphthalene Pipe not encountered. Top of water table encountered at approximately 5.5 ft bgs
1		
2		
3		
4		
5	Wet, dark brown/grey fine to coarse SAND and fine to coarse sub-angular to sub- round Gravel, some Slag, little Metal, trace Glass	
6	Bottom of Test Pit at 6 ft bgs	
7		
8		
9		

PROJECT NAME: Con Edison: Ludlow Street Works

PROJECT NUMBER:

446110

WEATHER

Partly cloudy, 60s, 5-10 mph E

November 7, 2012/0910 DATE/TIME START: DATE/TIME FINISH: November 7, 2012/1120

CONTRACTOR: ADT

INSPECTOR: Zohar Lavy TEST PIT ID:

TP-6

LOCATION:

Approximate L X W X D

= 29' X 4' X 6'

West of MTA Tracks, along Federal Street

PHOTOGRAPH

TP-6 looking South



TP-6 at 6 ft bgs. Water intrusion at approximately 5.5 ft bgs

TP-6 looking east



TP-6 Commencing excavation

	PARSONS TEST PIT RECORD	
PROJECT NAME:	Con Edison: Ludlow Street Works	TEST PIT ID: TP-1
PROJECT NUMBER:	446110	LOCATION:
WEATHER	Partly cloudy, 60s, 5-10 mph E	Approximate L X W X D
DATE/TIME START:	June 17, 2010/1230	= 29' X 4' X 8.5'
DATE/TIME FINISH:	June 17, 2010/1400	
CONTRACTOR:	ECI	Northwest portion of the site
INSPECTOR:	Zohar Lavy	
DEPTH (feet bgs)	FIELD IDENTIFICATION OF MATERIAL	COMMENTS
0	0-0.5' ASPHALT	Holder foundation wall not encountered. Holder bottom/floor encountered at ~ 8.5 ft
1	0.5-3' Dark brown fine to coarse SAND, some Gravel and Cobble, little Brick, trace Boulder and Wood	bgs.
2		
3	3-6' Dark brown medium to coarse SAND, some Gravel and Brick, trace Cobble	1
4		
5		
6	6-8.5' ASH and SLAG, little fine to coarse Sand, Cobble and Brick	
7		
8		
	Bottom of Test Pit at 8.5 ft bgs	
9		
10		
11		
12		

PROJECT NAME: Con Edison: Ludlow Street Works

446110 PROJECT NUMBER:

WEATHER

Partly cloudy, 60s, 5-10 mph E DATE/TIME START: June 17, 2010/1230

DATE/TIME FINISH: June 17, 2010/1400

CONTRACTOR: ECI

INSPECTOR: Zohar Lavy TEST PIT ID: TP-1

LOCATION:

Approximate L X W X D

= 29' X 4' X 8.5'

Northwest portion of the site

PHOTOGRAPH

TP-1looking west



TP-1 Close up looking east



TP-1 Scraped holder bottom at 8.5 ft bgs

PARSONS							
TEST	PIT	RECOR	D				

Г

	TEST PIT RECORD	
PROJECT NAME:	Con Edison: Ludlow Street Works	TEST PIT ID: TP-2
PROJECT NUMBER:	446110	LOCATION:
WEATHER	Clear, 70s, 0-5 mph East	Approximate L X W X D
DATE/TIME START:	June 18, 2010/0800	= 50' X 4' X 4.5'
DATE/TIME FINISH:	June 18, 2010/1215	Angled test-pit = 8' X 4' X 8.5'
CONTRACTOR:	ECI	
INSPECTOR:	Zohar Lavy	Northwest portion of the site, just south of junk pile
DEPTH (feet bgs)	FIELD IDENTIFICATION OF MATERIAL	
	0-4.5' CONSTRUCTION DEBRIS and brown fine to coarse SAND, some Brick; 0-4.5' (in Angled Test Pit) BRICK and brown fine to coarse SAND	Holder foundation wall not encountered in originalTP-2. Footings encountered at ~4.5 ft bgs in original
1		TP-2. Holder wall encountered from ~3.5 to 8.5 ft bgs in angled TP-2 (~18 ft east of TP-2 western starting point).
2		Holder foundation wall encountered again at ~ 40 ft east from TP-2 western starting point.
3		western starting point.
4	4.5-7.5' (in Angled Test Pit) BRICK and brown fine to coarse SAND	
5	4.5-7.5 (III Aligied Test Fit) Bittort and blown line to coarse SAIVE	
6		
7		
8	7.5-8.5' (in Angled Test Pit) ASH and SLAG, little fine to coarse Sand	Holder bottom/floor encountered at ~8.5 ft bgs
	Bottom of Test Pit at 8.5 ft bgs	

PROJECT NAME: Con Edison: Ludlow Street Works

PROJECT NUMBER:

446110

WEATHER

Clear, 70s, 0-5 mph East

DATE/TIME START: DATE/TIME FINISH: June 18, 2010/0800 June 18, 2010/1215

CONTRACTOR:

CONTRACTOR: <u>E</u>
INSPECTOR: Z

ECI Zohar Lavy TEST PIT ID: TP-2

LOCATION:

Approximate L X W X D

= 50' X 4' X 4.5'

Angled test-pit = 8' X 4' X 8.5'

Northwest portion of the site, just south of

junk pile

PHOTOGRAPH

TP-2 looking northwest



TP-2 looking west

TP-2 (angled) at ~ 8.5 ft bgs with holder foundation wall in foreground and footings in back ground



TP-2 eastward extension with exposed holder foundation wall and larger excavation in back ground

	PARSONS TEST PIT RECORD	
PROJECT NAME:	Con Edison: Ludlow Street Works	TEST PIT ID: TP-3
PROJECT NUMBER:	446110	
WEATHER	Partly cloudy, 60s, 5-10 mph east	LOCATION:
DATE/TIME START:	June 17, 2010/0800	Approximate L X W X D
DATE/TIME FINISH:	June 17, 2010/1130	= 11' X 4' X 9.5'
CONTRACTOR:	ECI	Central-Eastern portion of the site
INSPECTOR:	Zohar Lavy	
DEPTH (feet bgs)	FIELD IDENTIFICATION OF MATERIAL	COMMENTS
0	0-0.5' ASPHALT	Holder wall encountered at ~ 0.5 ft bgs and intact to 9.5 ft bgs
1	0.5-3' Brown fine to medium SAND, some Gravel, little sub-round Cobble, trace Urban fill, dry	
2		
3	3-6' Brown/red medium to coarse SAND, some angular and sub-angular Gravel, trace Cobble, dry	
4		
5		
6	6-9' Dark brown/red fine to coarse SAND, little sub-angular Gravel, little Concrete and Brick, trace Boulder, slightly moist	
7		
8		
9	9-9.5' Dark brown/black fine to coarse SAND, little sub-angular Gravel, moist	Holder bottom encountered at ~ 9.5 ft bgs
10	Bottom of Test Pit at 9.5 ft bgs	
11		
12		

PROJECT NAME: Con Edison: Ludlow Street Works

PROJECT NUMBER: 446110

WEATHER Partly cloudy, 60s, 5-10 mph east

DATE/TIME START: June 17, 2010/0800

DATE/TIME FINISH: June 17, 2010/1130

CONTRACTOR: ECI

INSPECTOR: Zohar Lavy

TEST PIT ID: TP-3

LOCATION:

Approximate L X W X D

= 11' X 4' X 9.5'

Central-eastern portion of the site

PHOTOGRAPH

TP-3 looking north



TP-3 looking north into excavation

TP-3 looking north TP-3 and holder wall exposed at ~0.5 ft bgs



TP-3 and holder wall exposed to 9.5 ft bgs

	PARSONS TEST PIT RECORD	
PROJECT NAME:	Con Edison: Ludlow Street Works	TEST PIT ID: TP-4A/B
PROJECT NUMBER:	446110	
WEATHER	Partly cloudy, 60s, 0-5 mph east	LOCATION:
DATE/TIME START:	June 14, 2010/1100	Approximate L X W X D
DATE/TIME FINISH:	June 14, 2010/1425	TP-4A = 12' X 4' X 3'
CONTRACTOR:	ECI	TP-4B = 10' X 4' X4.5'
INSPECTOR:	Zohar Lavy	Northern portion of the site
DEPTH (feet bgs)	FIELD IDENTIFICATION OF MATERIAL	COMMENTS
0	0-0.5' SALT	Gasholder wall not encountered. A brick structure with two ~12" diameter pipes
1	0.5-4.5' Brown fine to coarse SAND, some Gravel, some Brick, little Cobble, little Urban fill and Debris	running northeast/southwest was exposed. The holder wall was later uncovered at the surface ~8 ft northeast of TP-4B during salt moving operations. Additionally, two
2		open ~12" pipes were encountered ~ 2 ft north of TP-4B.
3		
4		
5	Bottom of Test Pit at 4.5 ft bgs	
6		
7		
8		
9		
10		
11		

PROJECT NAME: Con Edison: Ludlow Street Works

PROJECT NUMBER: 446110

WEATHER Partly cloudy, 60s, 0-5 mph east

DATE/TIME START: <u>June 14, 2010/1100</u>

DATE/TIME FINISH: June 14, 2010/1425

CONTRACTOR: ECI

INSPECTOR: Zohar Lavy

TEST PIT ID: TP-4A/B

LOCATION:

Approximate L X W X D

TP-4A = 12' X 4' X 3'

TP-4B = 10' X 4' X4.5'

Northern portion of the site

PHOTOGRAPH

TP-4A looking north



TP-4A and 4B looking southwest

TP-4A and brick structure to the north with 2 ~12" pipes running southwest from under brick structure



TP-4B in foreground and TP-4A in background with the square brick structure adjacent to both

	PARSONS TEST PIT RECORD	
PROJECT NAME:	Con Edison: Ludlow Street Works	TEST PIT ID: TP-5
PROJECT NUMBER:	446110	
WEATHER	Clear, 60s, 0-5 mph east	LOCATION:
DATE/TIME START:	June 15, 2010/0720	Approximate L X W X D
DATE/TIME FINISH:	June 15, 2010/1145	12' X 4' X 9'
CONTRACTOR:	ECI	
INSPECTOR:	Zohar Lavy	Northern portion of the site
DEPTH (feet bags)	FIELD IDENTIFICATION OF MATERIAL	COMMENTS
	0-5' Brown fine to coarse SAND, some Cobble, some Gravel, little Boulder, trace Brick, trace Urban debris, dry	Holder wall encountered at ~ 1.5 ft bgs. The bottom/floor of the holder was not encountered.
1		
2		
3		
4		
5	Brown fine to coarse SAND and COAL, little Gravel, some Cobble, dry	
6		
7	Brown/grey fine to coarse SAND and fine GRAVEL, little Gravel, trace Rubber, trace Metal	e Water encountered at ~ 9 ft bgs
8		
9	Bottom of Test Pit at ~ 9 ft bgs	
10		
11		
12		

PROJECT NAME: Con Edison: Ludlow Street Works

PROJECT NUMBER:

446110

WEATHER DATE/TIME START: Clear, 60s, 0-5 mph east

TE/TIME START: June 15, 2010/0720

DATE/TIME FINISH:

June 15, 2010/1145

CONTRACTOR: INSPECTOR: ECI Zohar Lavy TEST PIT ID: TP-5

LOCATION:

Approximate L X W X D

12' X 4' X 9'

Northern portion of the site

PHOTOGRAPH

TP-5 looking northeast



TP-5 looking southwest

TP-5 to 9 ft bgs with holder wall in the foreground



TP-5 to 9 ft bgs with holder wall reaching down to and passed water at 9 ft bgs

					PARSONS	BORING/WELL I	
Contracts	or: Aquifer Dril	lling & Tasting			DRILLING RECORD	Location Description:	Sheet 1 of 1
Driller:	Greg Rivera				PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Walkway immediately so	outh of Worselly Dron
Inspector	spector: Zohar Lavy			PROJECT NUMBER: 446110-04000	Building	duit of waverry Frop.	
Rig Type	Track Mour	nted Sonic Rig					
		WATER OBSERVAT	TIONS			Location Plan	
Water	DTW	DTW			Weather: Overcast, low 60s		
Level	~23' bgs						
Date	10-26-2012				Date/Time Start: 10-26-2012/1415	See Site Plan	
Time	1450	m 60 :			D (75' F' 1 10 27 2012/1455		
Meas. From	ft bgs - Soil cuttings	Top of Casing			Date/Time Finish: 10-26-2012/1455	4	
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.	511	(%)	(ppm)		JOHE MITTO	00
+2			()	, T			
+1							
0		Vac-tron	NA	NA	0-4" ASPHALT		
1		Vac-tron	NA	0.0	4"-5' Dry, fine to coarse SAND and DEBRIS, some fine to coarse sub-angular to sub-round Gravel,		
2		Vac-tron	NA NA		little Asphalt		
3		Vac-tron Vac-tron	NA NA				
5		vac-tion	INA		0-12" Dry, dark brown fine to coarse SAND, some fine to coarse sub-angular Gravel, trace Brick		
6	SB-9 (5-8)	Sonic	83	9.4	fragments		
7					12-30" Dry, browm, fine to medium SAND and angular to round fine to medium Gravel		
8					0-12" SCHIST	1	Very hard drilling at 8' bgs
9		Sonic	100	8.2	12-24" Moist, brown/orange fine to medium SAND and fine to coarse sub-angular to sub-round		
					GRAVEL, trace Silt	4	
10		_			0-24" Dry/moist, brown medium SAND, some angular to sub-round fine to medium Gravel, trace Porcelain & Shell fragments		
11		Sonic	100	5.8	24-60" Dry, orange fine to medium SAND, little coarse sub-angular Gravel, trace Silt		
13		Bonic	100	3.0	, , , , , , , , , , , , , , , , , , ,		
14		†					
15					0-24" Dry, orange fine to medium SAND, little coarse sub-angular Gravel, trace Silt	1	
16					24-50" Dry, orange/brown medium to coarse SAND, some angular to sub-angular fine to medium		
17		Sonic	100	3.5	Gravel		
18		_			50-60" Moist, orange/brown fine to coarse SAND, some medium to coarse sub-round Gravel, trace Silt		
19 20	 	 	+		0-24" Moist, orange/brown fine to coarse SAND, some medium to coarse sub-round Gravel, trace Silt, trace Ash and	-	
21	 	1			Plastic	'	
22	<u> </u>	Sonic	80	4.4	24-48" Moist/wet, brown fine to coarse SAND, some medium to coarse sub-round Gravel, trace Silt		
23]					
24							
25		1]	0-24" Moist/wet, brown fine to coarse SAND, some medium to coarse sub-round Gravel, little Silt		
26	-	Sonic	100	2.5	24-60" Moist, orange fine SAND and Silt, trace Clay and Cobble		
27 28		Sonic	100	2.5			
29	SB-9 (28-30)						
30			1		End of Boring at 30 ft bgs	1	
	SAMPLING MET	HOD			Hand cleared to 5 ft bgs; Sonic drilled from 5 to 30 ft bgs		,
	SS = SPLIT SPOON	ı					
	A = AUGER CUTT	INGS					
	C = CORED	HAMMED (DODG)					
	WH = WEIGHT OF	HAMMER (RODS)					

Track Mounted Sonic Rig						PARSONS	BORING/WELL II	D: SB-10
March Complete C	_					DRILLING RECORD		Sheet 1 of 1
Part						PROJECTNAME Consultated Editors Francis Ledling Const Weeks		V1
About 1	Driller:	Greg Rivera				PROJECT NAME: Consolidated Edison - Former Ludiow Street Works		
CROCKINNATER ORDERVATIONS CROSSINVATIONS CROSSINVAT	Inspector	: Zohar Lavy				PROJECT NUMBER: 446110-04000	-	,
Western DFW DFW DFW December Dec	Rig Type:							
Local Section 1925 1920				TONS	1		Location Plan	
Note 19.25 20.12			DTW			Weather: Overcast, low ous	_	
Date Description Descrip	Date					Date/Time Start: 10-25-2012/1125	See Site Plan	
SCHEMATIC COMMENTS SPT Rec. Comment Comment	Time							
Depth Sample D. Pipp			Top of Casing			Date/Time Finish: 10-25-2012/1240	1	
Sample LD. Chain Cympan Chain Cympan			SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
1								
1								
1			Vac-tron	N/A	0.0	0.6" Majet dark brown fine to coarse SAND and ORGANICS some fine to coarse Gravel trace Silt	_	
Vacation	1				0.0			
1			Vac-tron	NA	0.0			
1.5 1.5				_	0.0	Gravel 3- 5' Moist/dry brown, medium to fine SAND, some fine to coarse Gravel, trace Brick		
The content of the			Vac-tron	NA	0.5	0-52" Moist, dark brown fine to coarse SAND, some fine to coarse sub-angular to sub-round Gravel	1	
10			1		0.0			
9	7		Sonic	100	1.8	52-60" APHALT and WOOD, some fine to coarse Sand and fine to coarse Gravel		
10								
1	_			1		0-12" Dry brown/gray fine to coarse SAND, some fine to coarse Grayel, trace Wood and, Metal	-	
30.48* Mosts, brown medium to fine SAND and weathered schist COBBLE, little fine to coarse Gravel, trace Wood, Procedus, and Sill. 1.5 0.40* Mosts, brown, medium to fine SAND, some fine to coarse sub-angular to sub-round Gravel, little Asphals, Wood, trace Conceste 1.5 0.40* Mosts, brown, medium to fine SAND, some fine to coarse sub-angular to sub-round Gravel, little Asphals, Wood, trace Conceste 1.5 0.40* Mosts, brown, medium to fine SAND, some fine to coarse sub-angular to sub-round Gravel, little Asphals, Wood, trace Conceste 1.5 0.40* Mosts, brown, medium to fine SAND, some fine to coarse sub-angular to sub-round Gravel, little Asphals, Wood, trace Conceste 1.5 0.40* Mosts, black fine to coarse SAND and COAL, slight hydrocarbon odor 1.6 0.40* Mosts, orange-brown fine to coarse SAND, little angular to sub-nound medium to fine Gravel, trace Sill, trace Sill, 1.5 0.40* Mosts, daming, strong hydrocarbon odor, NAPL 0.40* Mosts, daming, strong			1					
Table	12		Sonic	80	1.2			
1.5 1.5 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.6 1.5 1.5 1.6 1.5 1.5 1.6 1.5						· · · · · · · · · · · · · · · · · · ·		
Asphalt, Wood, trace Concrete					1.5		_	
1					1.5			
9			Sonic	100				
1.3								
Asphalt, Wood, trace Concrete 222 23 24 25 26 26 27 26 27 27 27 27					1.3	0.16" Moist brown madium to fine SAND some fine to coarse sub-angular to sub-round Gravel little	-	
23					1.3			
17. 40.48* Moist, orange/brown fine to coarse SAND, little angular to sub-round medium to fine Gravel, trace Sit			Sonic	80	-			
100 100								
100 25.0 100 28.50 24.60° Wet, black fine to coarse SAND and angular to sub-angular medium to fine Gravel, some Coal, black staining, strong hydrocarbon odor, NAPL 22.54.0 024° Wet, black fine to coarse SAND and angular to sub-angular medium to fine Gravel, some Coal, black staining, strong hydrocarbon odor, NAPL 24.28° BRICK 28.34° CONCRETE 1200.0 34.60° Dry, orange-brown, fine to coarse SAND, little medium to fine sub-angular Gravel, trace Silt, 35.0 35.00 35.00 35.00 35.00 35.00 35.00 35.00 35.00 36.00° Dry, orange-brown, fine to coarse SAND, little medium to fine sub-angular Gravel, trace Silt, 35.00 36.00° Dry, orange-brown, fine to coarse SAND, little medium to fine sub-angular Gravel, trace Silt, 35.00 36.00° Dry, orange-brown, fine to coarse SAND, orange NAPL saturated, strong hydrocarbon odor and staining, strong hydrocarbon odor, NAPL blebs 36.00° Dry, orange-brown, fine to coarse SAND, orange NAPL saturated, strong hydrocarbon odor and staining, strong hydrocarbon odor and staining 36.00° Dry, orange-brown, fine to coarse GRAVEL and SAND, NAPL saturated, strong hydrocarbon odor and staining 36.00° Dry, orange fine SAND and SILT, trace olive Clay, sheen 36.00° Dry, orange fine SAND and SILT, trace olive Clay, sheen 36.00° Dry, orange fine SAND and SILT, trace olive Clay, sheen 36.00° Dry, orange from SAND, inthe Silt 36.00° Dry, orange from SAND, inthe Salt 36.00° Dry, orange from SAND, inthe Salt 36.00° Dry, orange from SAND, inthe SAND, inthe Salt 36.00° Dry, orange from SAND, inthe SAND, inthe Salt 36.00° Dry, orange from SAND, inthe Salt 36.00							-	
Sonic 100								
28			Sonic	100	835.0			
254 0 - 24" Wet, black fine to coarse SAND and angular to sub-angular medium to fine Gravel, some Coal. black staining, strong hydrocarbon odor, NAPL						black staining, strong hydrocarbon odor, NAPL		
Signature Sign	29							
32		SB-10 (30-32)			2254.0			
28-34" CONCRETE 1200.0 28-34" CONCRETE 1200.0 34-60" Dry, orange/brown, fine to coarse SAND, little medium to fine sub-angular Gravel, trace Silt, 34-60" Dry, orange/brown, fine to coarse SAND, little medium to fine sub-angular Gravel, trace Silt, Staining, strong hydrocarbon odor, NAPL blebs 350.0 20-60" Wet, black, medium to coarse SAND, orange NAPL saturated, strong hydrocarbon odor and staining 485.0 30" Wet, black fine to coarse GRAVEL and SAND, NAPL saturated, strong hydrocarbon odor 30-36" Wet, orange medium to coarse SAND 36-60" Wet, orange, fine SAND and SILT, trace olive Clay, sheen 44 45 58-10 (45-47) 46 58-10 (45-47) 47 50nic 100 48 50 50 50 50 50 50 50 5			Sonic	100				
1200.0 34-60° Dry, orange/brown, fine to coarse SAND, little medium to fine sub-angular Gravel, trace Silt,			Some	100		28-34" CONCRETE		
Sampling Methods Sonic S				ļ	1200.0	34-60" Dry, orange/brown, fine to coarse SAND, little medium to fine sub-angular Gravel, trace Silt,	1	
350	35				1150.0			
Sonic 100 Staining Staini	36			100	350.0			
39			Sonic	100				
40			-					
30-36" Wet, orange medium to coarse SAND 36-60" Wet, orange, fine SAND and SILT, trace olive Clay, sheen 43				+	485.0	0-30" Wet, black fine to coarse GRAVEL and SAND, NAPL saturated, strong hydrocarbon odor	1	
42 Sonic 100 30-60 Wet, orange, line SAND and SILT, trace olive Clay, sheen	41]			30-36" Wet, orange medium to coarse SAND		
44			Sonic	100	03.0	36-60" Wet, orange, fine SAND and SILT, trace olive Clay, sheen		
45			-					
46		CD 10 /45 15		1	22.0	0-12" Moist/wet, orange, fine SAND and SILT, trace olive Clay, sheen	†	
48	46	SB-10 (45-47)				12-24" Moist, orange/red fine SAND, little Silt		
49 End of Boring at 50 ft bgs			Sonic	100		24-60" Weathered SCHIST		
End of Boring at 50 ft bgs SAMPLING METHOD NAPL observed from 27 to 42.5 ft bgs SS = SPLIT SPOON A = AUGER CUTTINGS C = CORED End of Boring at 50 ft bgs NAPL observed from 27 to 42.5 ft bgs NAPL observed from 27 to 42.5 ft bgs Soring hand cleared to 5 ft bgs; Sonic drilled 5 to 50 ft bgs C = CORED			-					
SAMPLING METHOD NAPL observed from 27 to 42.5 ft bgs SS = SPLIT SPOON A = AUGER CUTTINGS C = CORED NAPL observed from 27 to 42.5 ft bgs Boring hand cleared to 5 ft bgs; Sonic drilled 5 to 50 ft bgs C = CORED			<u>I</u>	1	I	End of Boring at 50 ft bgs	1	
A = AUGER CUTTINGS C = CORED		SAMPLING MET	HOD			NAPL observed from 27 to 42.5 ft bgs	•	•
C = CORED						Boring hand cleared to 5 ft bgs; Sonic drilled 5 to 50 ft bgs		
			IIVOS					
			HAMMER (RODS)					

					PARSONS	BORING/WELL II	
Contract	ontractor: Aquifer Drilling & Testing			DRILLING RECORD	L di D i . di	Sheet 1 of 1	
					BROJECT NAME: Consolidated Edison: Former Ludlow Street Works	Location Description:	
Driller:	Greg Rivera	<u> </u>			PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	On Federal Street immed	iately west of the
Inspector	spector: Matt Bruno			PROJECT NUMBER: 446110-04000	MTA tracks and south of	the Sugar Refinery	
Rig Type		nted Sonic Rig			-		
***		WATER OBSERVAT	TIONS		D - 1 1 10	Location Plan	
Water Level	DTW ~5' bgs	DTW			Weather: Partly cloudy, 40s	-	
Date	11-06-2012				Date/Time Start: 11-06-2012/1120	See Site Plan	
Time	1130				Date Time Start.	- See Site 1 km	
Meas.	ft bgs -	Top of Casing			Date/Time Finish: 11-06-2012/1150		
From	Soil cuttings						
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.		(%)	(ppm)			
+2			-				
+1 0	1	Vac-tron	NA	0.0	0-12" Wet, black fine to medium SAND and sub-angular to sub-round GRAVEL	4	
1		Vac-tron	NA NA	0.0	12-5' Dry/moist, brown fine to coarse SAND and weathered Schist COBBLES/BOULDERS, some fine		
2		Vac-tron	NA	0.0	to coarse sub-angular Gravel, trace Metal and Slag		
3		Vac-tron	NA				
4		Vac-tron	NA				
5	SB-11 (5-5.5)			1.5	0-24" Wet, black/dark brown, angular GRAVEL and CINDERS, little Silt	1	
6				1.9	24-36" Wet, black SILT, little angular Cobble, trace Coal fragments		
7		Sonic	60				
8		4					
9				3.7	0-24" Wet, black, soft fine SAND, little Silt and sub-angular Gravel, trace Plastic fragments, slight	4	
11				5.7	hydrocarbon odor		
12		Sonic	80	4.2	24-48" Wet, red/brown medium SAND, trace Silt		
13		1					
14							
15				12.7	0-15" Wet, black/dark brown, sub-angular GRAVEL, little Silt		
16				15.2	15-30" Wet, black, medium SAND, little sub-angular Gravel, little Silt		
17		Sonic	50				
18 19	1	-					
20	+	 	+	20.2	0-12" Moist, red/brown, soft, medium SAND, some Silt	1	
21		1		27.3	12-48" Wet, red/brown, medium dense, SILT, little Mica Schist		
22		Sonic	80				
23]					
24						1	
25		1		0.0	0-6" Moist, red/brown SILT, lense of white/tan micaceous Sandstone, trace Mica Schist		
26		G	90	0.0	6-48" Dry, red/brown very dense SILT, trace sub-round Gravel		
27 28	1	Sonic	80				
28	SB-11 (29.5-30)	4					
30	22 11 (2).0 30,	1	1	ı	End of boring at 30 ft bgs	1	
30	SAMPLING MET	HOD			Hand cleared to 5 ft bgs; Sonic drilled from 5 to 30 ft bgs		
	SS = SPLIT SPOON						
	A = AUGER CUTT	INGS					
	C = CORED						
	WH = WEIGHT OF	HAMMER (RODS)					

					PARSONS	BORING/WELL II	D: SB-12
					DRILLING RECORD		Sheet 1 of 1
	Aquifer Dril				PROJECTIVANE C. III. IEE. E. J. H. C. W. I	Location Description:	
Driller:	Greg Rivera	· 			PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Approx. 25' west and in l	ine w/ DPW yard
Inspector					PROJECT NUMBER: 446110-04000	concrete retaining wall.	
Rig Type:		ited Sonic Rig					
Water	DTW	VATER OBSERVAT DTW	IONS	I	Weather: Overcast, low 60s	Location Plan	
Level	~22.5' bgs	51,,			Transfer and the second	1	
Date	11-05-2012				Date/Time Start: 11-05-2012/0950	See Site Plan	
Time Meas.	1450 ft bas	Top of Casing			Date/Time Finish: 11-05-2012/1100		
From	ft bgs - Soil cuttings	Top of Casing			Date/Time Philsit. 11-05-2012/1100	-	
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.		(%)	(ppm)			
+2			+				
0		Vac-tron	NA	0.0	0-1.5' Wet, dark brown, fine to medium SAND, some Debris, little Cobble, fine to coarse sub-angular to	†	
1		Vac-tron	NA		round Gravel, trace Brick and Silt		
3		Vac-tron	NA	0.0	1.5-5' Moist, dark brown, fine to medium SAND, some Cobble, little Brick, trace Silt		
4		Vac-tron Vac-tron	NA NA				
5		7.00			0-14" CONCRETE	1	
6		g .	50	- 1	14-18" BRICK		
7 8		Sonic	50	5.4	18-30" Moist, orange/brown fine to medium SAND, some fine to medium sub-angular Gravel		
9							
10				135.0	0-20" Moist, orange/grey fine to medium SAND, some fine to medium sub-angular Gravel, hydrocarbon		
11				1000	odor 20-50" Moist/wet, orange/grey fine to medium SAND, some fine to medium sub-angular Gravel, trace		
12	SB-12 (12-14)	Sonic	100	1977.0	Cobble, NAPL saturated, hydrocarbon and napthalene odor and staining		
14		1		721.0	50-60" Moist, tan/orange SILT, some fine Sand, little olive Clay, little fine to medium Sand lenses,		
				721.0	hydrocarbon odor 0-54" Moist, orange SILT and fine SAND, slight hydrocarbon odor	-	
15 16		•			54-60" Moist, orange SILT and fine SAND, little medium Sand, slight hydrocarbon odor		
17		Sonic	100	65.5			
18							
19 20					0-30" Moist, orange/brown SILT and fine SAND, little tan Clay, little fine to medium Sand lenses,	-	
21		1			hydrocarbon odor		
22		Sonic	50	73.7			
23							
24 25			+		0-30" Moist, orange/brown SILT and fine SAND, little tan Clay, little fine to medium Sand lenses,	1	
26]			hydrocarbon odor		
27		Comio	100	41.6	30-36" Moist, orange/brown SILT and fine SAND, little tan Clay, little fine to medium Sand lenses,		
28		Sonic	100	41.6	hydrocarbon odor, sheen 36-52" Moist, orange/brown SILT and fine SAND, little tan Clay, little fine to medium Sand lenses,		
29					hydrocarbon odor		
20				10:-	52-60" Moist, orange medium SAND, slight hydrocarbon odor	4	
30				121.0 68.4	0-32" Wet, orange/brown medium to coarse SAND, hydrocarbon odor and slight staining 32-60" Wet, orange/brown medium to coarse SAND, hydrocarbon odor, trace fine sub-angular to round		
31		Sonic	100	06.4	Gravel		
33		j					
34						1	
35 36					0-24" Wet, orange/brown medium to coarse SAND, hydrocarbon odor, little fine sub-angular to round Gravel trace fine sub-angular to round Gravel		
37	SB-12 (36-38)	Sonic	75	22.9	24-36" Weathered white SCHIST, NAPL within fractures		
38							
39					End of boring at 39 ft bgs	1	
	SAMPLING MET SS = SPLIT SPOON				NAPL observed from 36 - 38 ft bgs Hand cleared to 5 ft bgs; Sonic drilled from 5 to 39 ft bgs		
	A = AUGER CUTT						
	C = CORED						
	WH = WEIGHT OF	HAMMER (RODS)					

					PARSONS	BORING/WELL II	
_					DRILLING RECORD		Sheet 1 of 1
	r: Aquifer Drill					Location Description:	
Driller:	Greg Rivera				PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Approximately 70' north concrete retaining wall an	
Inspector:	Zohar Lavy				PROJECT NUMBER: 446110-04000	west of Waverly Properti	
Rig Type:	Track Moun	ted Sonic Rig				, ,	•
		VATER OBSERVAT	IONS			Location Plan	
Water	DTW	DTW			Weather: Partly cloudy, 40s		
Level	~22' bgs 10-26-2012				D (T)	G G' DI	
Date Time	1450				Date/Time Start: 11-05-2012/1420	See Site Plan	
	ft bgs -	Top of Casing			Date/Time Finish: 11-05-2012/1505		
From	Soil cuttings						
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth +2	Sample I.D.		(%)	(ppm)			
+1							
0		Vac-tron	NA	0.0	0-12" Dry, dark brown, fine to coarse SAND, some Organics (rootlets), little Gravel and Cobble	1	
1		Vac-tron	NA	0.0	12-5' Dry brown/orange fine to coarse SAND and COBBLE, some Brick, little Concrete fragments		
3		Vac-tron Vac-tron	NA NA				
4		Vac-tron Vac-tron	NA NA				
5			- '		0-12" Weathered gneissic SCHIST boulder	1	
6					12-30" Dry, dark brown, fine to medium SAND and fine to coarse angular to sub-round GRAVEL		
7		Sonic	50	1.3			
9							
10					Dry, dark brown, fine to medium SAND and fine to coarse angular to sub-round GRAVEL some	1	
11			400		gneissic Schist Cobbles and Organics		
12		Sonic	100	1.7			
14							
15					0-10" Dry, dark brown, fine to medium SAND and fine to coarse angular to sub-round GRAVEL some	1	
16					gneissic Schist Cobbles and Organics, trace Silt		
17 18		Sonic	100	2.1	10-48" SCHIST Boulder 48-60" Moist, dark brown/black fine to medium SAND, some fine to coarse sub-angular to sub-round		
19					Gravel		
20					0-20" Dry, gray fine to coarse SAND, some sub-angular to sub-round Gravel, little Schist Cobble	1	
21			400		20-40" CONCRETE		
22 23		Sonic	100	2.3	40-60" Moist, black, fine to medium SAND and BRICK, some fine to medium sub-angular Gravel, trace Glass fragments		
24							
25					0-12" Moist, orange/brown fine to medium SAND, little Silt, trace sub-angular to round fine to coarse		
26		G :-	100	1.5	Gravel 12-60" Dry, orange/brown medium to coarse SAND, little coarse angular Gravel, trace Schist Cobble		
27 28		Sonic	100	1.3	12-00 Dry, orange/brown medium to coarse SAND, inthe coarse angular Graver, trace Schist Cobble		
29							
30					0-10" Dry, orange/brown medium to coarse SAND, little coarse angular Gravel, trace Schist Cobble		
31		Somia.	100	2.1	10-60" Moist orange/brown fine to medium SAND, some Silt, trace course round gravel		
32		Sonic	100	2.1			
34							
35					0-36": Moist, orange fine SAND and SILT		
36 37		Sonic	100	66.2	36-60" Moist, orange/gray fine SAND, slight hydrocarbon odor		
38		Some	100	00.2			
39							
40				73.7	0-24" Moist, grey/black fine to medium SAND, some angular Gravel, little Schist, trace Silt, slight		
41 42					hydrocarbon odor 24-40" Moist/wet, orange fine to medium SAND and weathered gneissic SCHIST COBBLE		
43				85.6	40-50" Moist/wet, orange fine to medium SAND and weathered gnessic SCHIST COBBLE, NAPL		
44		Sonic	46		blebs and sheens		Rock encountered at
45							approximate;y 48 ft bgs
46 47							
48	SB-13 (47-49)						
49					End of boring at 49 ft bgs	1	
	SAMPLING METI				NAPL observed from 47 - 49 ft bgs		
	SS = SPLIT SPOON A = AUGER CUTTI				Hand cleared to 5 ft bgs; Sonic drilled from 5 to 49 ft bgs		
	A = AUGER CUTTI C = CORED						
	WH = WEIGHT OF	HAMMER (RODS)					

					PARSONS	BORING/WELL II	
.	4 :c 15 :T				DRILLING RECORD		Sheet 1 of 1
	r: Aquifer Dril					Location Description:	
Driller:	Jeremey Me	yers			PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Approximately 100' north	
Inspector	Zohar Lavy				PROJECT NUMBER: 446110-04000	concrete retaining wall ar west of Waverly Properti	
Rig Type:		ted Sonic Rig					,,.
Kig Type.		VATER OBSERVAT	IONS			Location Plan	
Water	DTW	DTW			Weather: Rain, 40s, 15-20 mph East		
Level	~36' bgs						
Date	12-21-2012				Date/Time Start: 12-21 -2012/0925	See Site Plan	
Time	1050	m 40 i			D . Th. 10.01.0010/1170		
	ft bgs - Soil cuttings	Top of Casing			Date/Time Finish: 12-21-2012/1150	_	
From Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.	511	(%)	(ppm)	THE DESCRIPTION OF MATERIAL	SCHEMITTE	COMMENTS
+2							
+1							
0		Vactron	NA	0.0	0-5' Moist, brown, fine to coarse SAND, some Boulder, little Concrete, little fine to coarse Gravel, trace	,	
1		Vactron Vactron	NA NA		Plastic		
3		Vactron	NA NA				
4		Vactron	NA				
5			1	0.0	0-12" Wet, dark brown fine to medium SAND and COBBLE	1	
6				0.0	12-20" Wet, sub-angular to angular GRAVEL		
7		Sonic	67	0.0	20-40" Moist, brown fine to medium SAND, some Marble (debris), little Silt		
8							
10				0.1	0-30" Wet, brown, fine to medium SAND, some fine to coarse angular to sub-round Gravel, trace Rope	-	
11				0.1	and Wood		
12		Sonic	80	0.1	30-48" Dry, brown/grey fine to medium SAND and SHALE (debris)		
13							
14							
15				0.1	0-20" Moist, brown/grey fine to medium SAND and SHALE (debris), some gneissic Schist, little		
16 17		Sonic	33		Asphalt		
18		Some	33				
19							
20				0.0	Moist, dark grey, fine to medium SAND and angular coarse GRAVEL, some gneissic Schist and		
21					Concrete		
22		Sonic	27				
23 24							
25			+	0.2	0-12" Moist, dark grey, fine to medium SAND and angular coarse GRAVEL, some gneissic Schist and	=	
26					Concrete		
27		Sonic	83	0.2	12-40" Black gneissic SCHIST		
28				0.2	40-50" Dry, red/brown medium SAND		
29				0.2	O COULD C. CANTO For I. I. I.C. I. C. C.	_	
30				0.2	0-20" Dry, orange, fine SAND, little sub-round Gravel, trace Silt 20-60" Dry, orange/brown, medium to coarse SAND, little fine to coarse sub-angular to sub-round		
32		Sonic	100	0.2	Gravel		
33							
34							
35				0.0	0-14" Moist, orange, fine SAND, some Silt, little fine to coarse sub-angular to sub-round Gravel		
36 37		Sonic	100	-	14-42" Moist-wet, orange, fine SAND and SILT, trace Clay 42-60" Moist-wet, dark orange, fine SAND, little Silt, trace Mica		
38		Some	100		142-00 MOISI-wet, dark orange, the SAND, little Silt, trace Mica		
39	SB-14 (38-40)						
40			1	0.2	Moist-wet, dark orange, fine SAND, little Silt, trace Mica	1	
41	_						[
42		Sonic	80				
43				-			
44			+	0.0	0-14" Moist-wet, dark orange, fine SAND, little Silt, trace Mica	4	Hard Drilling from 44-45 ft bgs
45 46	SB-14 (45-47)	Sonic	60	0.0	14-33" Moist, orange/brown, fine to medium SAND		Very hard drilling from 47-
47					33-36" SCHIST		48 ft bgs
48			•		End of boring at 48 ft bgs		
	SAMPLING METI				Hand cleared to 5 ft bgs; Sonic drilled from 5 to 48 ft bgs		
	SS = SPLIT SPOON						
	A = AUGER CUTT C = CORED	INGS					
		HAMMER (RODS)					
							4

					PARSONS	BORING/WELL I	D: SB-16
_					DRILLING RECORD		Sheet 1 of 1
	Aquifer Dril				PROJECTNAME Constituted Filters France Ledlers Const Weeks	Location Description:	f.Vl DDW
Driller:	Jeremey Me	yers			PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Approximately 40' north concrete retaining wall a	
Inspector	: Zohar Lavy				PROJECT NUMBER: 446110-04000	west of Waverly Propert	
Rig Type:	Track Moun	ted Sonic Rig				<u> </u>	
		VATER OBSERVAT	TIONS			Location Plan	
Water Level	DTW ~30' bgs	DTW			Weather: Partly cloudy, mid 40s, 0-5 mph North	4	
Date	~30 bgs 12-20-2012				Date/Time Start: 12-20-2012/1035	See Site Plan	
Time	1610				Ditto Time Suita	- See Sile 1 illin	
Meas.	ft bgs -	Top of Casing			Date/Time Finish: 12-20-2012/1710		
From	Soil cuttings	CDT	- P	DID	EIGH D IDENTIFICATION OF MATERIAL	COMPAND	COMPARING
Sample Depth	Location/ Sample I.D.	SPT	Rec.	PID (nnm)	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
+2	Sample 1.D.		(70)	(ррш)		†	
+1							
0		Vactron	NA	0.0	0-12" Moist, dark brown, fine to medium SAND and fine to coarse GRAVEL, some Organics (roots)		
2		Vactron Vactron	(%) (ppm) NA 0.0 0-12" Moist, dark brown, fine to medium SAND and fine to coarse GRAVEL, some Organics (roots) NA 0.0 12"-5' Moist, brown fine to coarse SAND, some Boulder, little Gravel NA NA 0.0 10" Moist, brown/grey fine to medium SAND, some Concrete 10-48" Dry, brown, fine to coarse SAND, some fine to coarse angular to sub-round Gravel, little 100 0.8 24-50" Dry, brown, fine to medium SAND and weathered gneissic SCHIST, some Concrete 100 0.8 24-50" Dry, brown, fine to medium SAND and weathered gneissic SCHIST, some Concrete 100 0.12" Moist, dark brown, fine to medium SAND and weathered gneissic SCHIST, some Concrete 100 0.8 24-50" Dry, brown, fine to medium SAND and weathered gneissic SCHIST, some Concrete 100 0.12" Moist, dark brown, fine to medium SAND, little Silt, little round to angular Gravel, trace 100 0.30" Moist, fine to coarse sub-angular GRAVEL and SCHIST, little fine to coarse brown Sand				
3		Vactron					
4		Vactron					
5							
6		Sonic	80	1.1	10-48" Dry, brown, fine to coarse SAND, some fine to coarse angular to sub-round Gravel, little		
7 8		Some	80	1.1			
9							
10						,]	
11		Conto	100	0.0			
12		Sonic	100	0.8			
14					50-00 Dry, of own, fine to medium SAND and weathered gheissic SCIIIS1, some Concrete		
15					0-12" Moist, dark brown, fine to medium SAND, little Silt, little round to angular Gravel, trace	1	
16					12-24" SCHIST		
17		Sonic	50	1.4			Very hard drilling from 17-
18 19		-					
20					0-30" Moist, fine to coarse sub-angular GRAVEL and SCHIST, little fine to coarse brown Sand	1	
21					30-60" BRICK		
22		Sonic	100	5.7			
23 24							
25					No Recovery	-	
26							Very hard drilling from 25-
27		Sonic	0	NA			30 ft bgs. Casing advanced to 30 ft bgs
28 29							
30					Wet, BRICK and fine to coarse angular to sub-angular GRAVEL, little fine to medium SAND, NAPL	4	
31	SB-16 (30-33)				blebs, moderate hydrocarbon odor, sheens		
32		Sonic	80	1406.0			
33							
34 35			-	-	0-14"Wet, BRICK and fine to coarse angular to sub-angular GRAVEL, little fine to medium SAND,	4	
36		1			little Silt, NAPL blebs, slight hydrocarbon odor, sheens		
37		Sonic	60	370.0	14-24" CONCRETE		
38					24-36" Wet, dark brown/black, medium to coarse SAND, slight hydrocarbon odor and staining		
39 40			+	1	0-40" Moist, orange/brown fine to medium SAND, slight hydrocarbon odor	4	
40		1			40-60" Moist, brown/grey, fine SAND, little Silt		
42		Sonic	100	0.9			
43]					
44			1		O 2011 Maries Incomplete Control City	4	
45 46	SB-16 (45-47)	Sonic	100	0.7	0-20" Moist, brown/grey, fine SAND, some Silt 20-24" SCHIST		
46		l	1		End of boring at 47 ft bgs	1	
· ·	SAMPLING MET	HOD			NAPL encountered from 30-36 ft bgs		
	SS = SPLIT SPOON				Hand cleared to 5 ft bgs; Sonic drilled from 5 to 47 ft bgs		
	A = AUGER CUTT C = CORED	INGS					
		HAMMER (RODS)					
		(/					

					PARSONS	BORING/WELL I	
G	4 10 TO 11				DRILLING RECORD		Sheet 1 of 1
Contracto		ling & Testing				Location Description:	
Driller:	Jeremey Me	yers			PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Approximately 125' nort concrete retaining wall a	
Inspector	Zohar Lavy				PROJECT NUMBER: 446110-04000	west of Waverly Propert	
Rig Type:	Track Moun	ted Sonic Rig					
rug Typer		VATER OBSERVAT	TONS			Location Plan	
Water	DTW	DTW			Weather: Partly cloudy, mid 40s, 0-5 mph North		
Level	~40' bgs						
Date	12-21-2012				Date/Time Start: 12-21-2012/1250	See Site Plan	
Time Meas.	1520 ft bgs -	Top of Casing	-		Date/Time Finish: 12-21-2012/1530		
From	Soil cuttings	Top of Casing			Date Time Finish. 12-21-2012/1330	-	
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.		(%)	(ppm)			
+2							
+1		***	27.4	0.0	O SIM I S S S S S S S S S S S S S S S S S	4	
1			13 ft				
2							
3		Vactron NA 0.0 Vactron NA Vactron NA Vactron NA Vactron NA O-12" Wet, grey/brown, fine to medium SAND, some angular to sub-round fine to coarse Gravel 12-20" BOULDER Sonic 100 0.0 20-60" Dry, dark brown, fine to coarse SAND and angular to sub-round fine to coarse GRAVEL, little Brick Very hard drilling fr 13 ft bgs. Casi advanced to 13 ft Gneissic SCHIST and CONCRETE, trace Wood Very hard drilling fr					
4		Vactron	NA				
5							
<u>6</u> 7		Conic	100	0.0			
8		Some	100	0.0			
9							
10						1	Very hard drilling from 10-
11							13 ft bgs. Casing
12					a L L agyram L agyram wy L	4	advanced to 13 ft bgs
13 14					Gneissic SCHIST and CONCRETE, trace Wood		
15		Sonic	50	0.0			
16							
17							Very hard drilling from 18-
18							20 ft bgs. Casing
19					Control SCHIET and CONCEPTE Annual World	4	advanced to 20 it bgs
20					Gheissic SCHIST and CONCRETE, trace wood		
22		Sonic	67	0.0			
23							
24							
25					Gneissic SCHIST and CONCRETE		
26 27		Sonic	33	0.0			
28		Bonie	33	0.0			
29							
30					0-34" Gneissic SCHIST and CONCRETE]	
31		g :	100	0.1	34-60" Dry, orange/brown, medium to coarse SAND and fine to medium sub-angular to sub-round		
32 33		Sonic	100	0.1	Gravel		
34	SB-17 (33-35)						
35			1		No Recovery	1	
36							
37		Sonic	0	NA			
38							
39 40			1	1	Moist/wet, brown/grey medium to coarse SAND, some fine to medium sub-angular to sub-round	1	
41					Gravel, trace Silt		
42		Sonic	20	0.0			
43							
44							
45 46	SB-17 (45-47)				0-24" Moist, brown/grey medium to coarse SAND, some fine to medium sub-angular to sub-round Gravel, trace Silt		Very hard drilling from 47
46		Sonic	50	0.0	24-30" Moist, brown/grey medium to coarse SAND, some fine to medium sub-angular to sub-round		48 ft bgs
7/					Gravel, little weathered Schist trace Silt		
48		-			End of boring at 48 ft bgs	1	
	SAMPLING METI				Hand cleared to 5 ft bgs; Sonic drilled from 5 to 48 ft bgs		
	SS = SPLIT SPOON						
	A = AUGER CUTTI C = CORED	NGS					
	WH = WEIGHT OF	HAMMER (RODS)					

					PARSONS	BORING/WELL ID	
					DRILLING RECORD		Sheet 1 of 1
Contracto		lling & Testing				Location Description:	
Driller:	Jeremey Me	eyers			PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Southeast portion of the Y	Onkers DPW vard
Inspector:				PROJECT NUMBER: 446110-04000			
Rig Type:						 	
		WATER OBSERVAT	ΓIONS			Location Plan	
Water	DTW	DTW			Weather: Partly cloudy, high 40s		
Level		6.87					
Date		12-26-12			Date/Time Start: 12-19-2012/1340	See Site Plan	
ime		0930					
Aeas.		Top of Casing			Date/Time Finish: 12-19-2012/1410		
rom							
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.		(%)	(ppm)			
+2							
+1							Locking J-plug on casing
0		Vac-tron	NA	0.0	0-12" Moist, black, fine to coarse SAND and ASPHALT	1 и I и	Flushmount manhole
1		Vac-tron	NA	0.0	12"-5' Moist, dark brown, fine to coarse SAND, some Brick, little Cobble, trace Metal, trace Silt		
2		Vac-tron	NA				Bentonite (0-2')
3		Vac-tron	NA				<u> </u>
4		Vac-tron	NA				2" PVC riser (0.5-4')
5				NA	Large piece of Metal		
6		1		- 1111	zambo proce or moral		
7		Sonic	13				
8							
9							2" ID PVC well screen
10					No Recovery	1 (00 1 000	0.002" slot (4-19')
11						3 3 3 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	
12		Sonic	0	NA			#2 silica (2-19')
13		1					"2 sinca (2 1))
14							
15					No Recovery	- 1 (33) - 1 (33)	
16					Two receivery		
17		Sonic	0	NA			
18		Some		11/1			PVC end cap 19'
19		1				Estat Basel	. ve end cap 17
20		1	+	 	0-30" Wet, dark brown/orange, fine to coarse SAND, little Silt, little fine to coarse sub-angular to sub-	4	
21		1			round Gravel		
22		Sonic	100	10.7	30-60" Moist, orange, fine SAND and SILT, some fine sub-angular to sub-round Gravel, trace Clay		
23		Some	100	10.7	50 to 1400st, orange, the SALVE and SEE1, some time sub-angular to sub-round Oravel, trace Clay		
		4					
24		I		<u> </u>	Fig. 1. Charles at 25th at	4	
25		won			End of boring at 25' bgs		
	SAMPLING MET				Hand cleared to 5' bgs, sonic drilled from 5' to 25' bgs.		
	SS = SPLIT SPOON						
	A = AUGER CUTT	INGS					
	C = CORED						
	WH = WEIGHT OF	HAMMER (RODS)					

					PARSONS DRILLING RECORD	BORING/WELL ID	: MW-5 (alt.) Sheet 1 of 1
Contracto	r: Aquifer Dril	ling & Testing				Location Description:	
Driller:	Jeremey Me				PROJECT NAME: Consolidated Edison - Former Ludlow Street Works		
Inspector		*			PROJECT NUMBER: 446110-04000	Northwest portion of the	Yonkers DPW yard
Rig Type:		ted Sonic Rig			-		
8 71		VATER OBSERVAT	TIONS			Location Plan	
Water	DTW	DTW			Weather: Partly cloudy, high 40s		
Level	~10 ft bgs	11.12			· · · ·	1	
Date	12-19-12	12-26-12			Date/Time Start: 12-19-12/ 1120	See Site Plan	
Time	1140	1315					
Meas.	ft bgs -	Top of Casing			Date/Time Finish: 12-19-12/ 1200		
From	Soil cuttings	Top of Cusing			Date I III I I I I I I I I I I I I I I I I	=	
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.	511	(%)	(ppm)	THE DIDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
+2	Sample 1.D.		(70)	(ppin)			
+1							Locking J-plug on casing
0		Vac-tron	NA	0.0	0-5' Moist, dark brown fine to coarse SAND, some Cobble, little fine to coarse sub-angular to round		Flushmount manhole
1		Vac-tron Vac-tron	NA NA	0.0	Gravel, little Brick, trace Metal		Fiushmount mannoie
				<u> </u>	Glaver, fittle Blick, flace Metal	1 20 00	D
2		Vac-tron	NA	1			Bentonite (1-5')
3		Vac-tron	NA				
4		Vac-tron	NA				
5				0.3	Moist, brown fine to coarse SAND, some sub-angular to round fine to medium Gravel, trace Fabric		2" PVC riser (0.5-7')
6						38-83	
7		Sonic	80			38_833	
8						3-8	
9						- 000 - 000	2" ID PVC well screen
10				2.0	0-36" Wet, orange/brown fine SAND, some Silt, trace Clay	3 3	0.002" slot (7-22')
11			400		36-48" Wet, orange/brown fine SAND and WOOD, little Silt	38-83	
12		Sonic	100	327.0	48-60" Black stained WOOD, strong Cresote odor		#2 silica (5-22')
13							
14						88	
15					0-38" Moist, orange SILT, some olive Clay, little fine Sand, trace round fine to medium Gravel	3 <u> </u> 3	
16					38-60" Moist, orange/brown medium to coarse SAND	[3 <u> _</u> 8 .	
17		Sonic	100	2.5			
18							
19				ļ			
20					Moist, orange/brown medium to coarse SAND, trace fine sub-round to round Gravel	<u> </u>	
21							PVC end cap 22'
22		Sonic	100	2.5			
23				1			
24						_	
25					End of boring at 25' bgs		
	SAMPLING METI	HOD			Hand cleared to 5' bgs, sonic drilled from 5' to 25' bgs.		
	SS = SPLIT SPOON						
	A = AUGER CUTTI	NGS					
	C = CORED						
	WH = WEIGHT OF	HAMMER (RODS)					

					PARSONS	BORING/WELL ID	
Contracto	or: Aquifer Drill	ing & Testing			DRILLING RECORD	Location Description:	Sheet 1 of 1
Driller:	Greg Rivera				PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	1	DPW yard concrete retaining
Inspector	Zohar Lavy				PROJECT NUMBER: 446110-04000	wall, 70' W of Wavery Pro	
Rig Type:		ed Sonic Rig					
	GROUNDW	ATER OBSERVAT	IONS			Location Plan	
Water Level	DTW	DTW 38.69'			Weather: Overcast, low 60s	-	
Date		11-09-2012/0745			Date/Time Start: 10-26-2012/0810	See Site Plan	
Time Meas.	ft bgs -	Top of Casing			Date/Time Finish: 10-26-2012/0920		
From	Split Spoon				-	1	
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
+2	Sample 1.D.		(70)	(ррш)			Locking J-plug on
+1		V 4	NIA	0.0	O C" Maint deal beauty for the second CANID and ODC ANICS are second for Court for Silver de	n 1	inner wall
0		Vac-tron Vac-tron	NA NA	0.0	0-6" Moist, dark brown fine to coarse SAND and ORGANICS, some fine to coarse Gravel, trace Silt and Asphalt		Flush Mount Well Cover and Concrete
2		Vac-tron	NA	0.0	6"-3' Moist, dark brown fine to coarse SAND and ORGANICS, some Cobble, little fine to coarse Gravel 3'-5' Moist/dry brown, medium to fine SAND, some fine to coarse Gravel, trace Brick		
4		Vac-tron Vac-tron	NA NA	0.0	3-3 Moisterly brown, medium to time SAND, some fine to coarse Graver, trace Brick		Grout (0-24')
5				5.6	0-36" Dry, brown, fine to coarse SAND, some sub-angular to round fine to coarse Gravel, trace Wood		
<u>6</u> 7		Sonic		4.2	and Cobble 36-48" Dry, brown, fine to coarse SAND, some sub-angular to round fine to coarse Gravel, littleWood,		
8		Some		4.2	trace Cobble		2-inch ID PVC Riser
9							(0.5-28')
10 11					0-20" Wet, brown/gray fine to medium SAND, little Silt, trace Wood & Concrete		
12					20-36" Gneiss BOULDER		
13							
15		Sonic	30	20.7			
16 17		Some	30	20.7			
18							
19							
20					0-24" Wet, brown/gray fine to coarse SAND, some fine to coarse angular to sub-angular Gravel, trace Silt		
22					and Wood		# 2 Sand (21'-33')
23		Sonic	40	14.7			
25							Bentonite Chips (24-26')
26 27				7.1	0-24" Moist, brown/gray fine to coarse SAND, some fine to coarse angular to sub-angular Gravel, trace Silt and Wood		_
28		Sonic	80		24-36" Dry, black, medium to fine SAND and fine to coarse angular to sub-round GRAVEL, some		
29				11.6	Concrete, trace Coal fragments 36-48" Dry, orange medium to fine SAND, trace angular to sub-angular fine Gravel		
30				5.2	0-30" Moist, orange/brown fine to coarse SAND and angular to round GRAVEL	% %	
32				11.5	30-60" Wet, orange medium to fine SAND, trace Silt		
33 34		Sonic	100				# 2 silica (26-48')
35							(23 13)
36 37	MW-7 (36-38)				0-56" Moist, orange fine SAND and SILT, some medium - fine Sand veins transmitting NAPL, striated tan Clay laminations, strong hydrocarbon odor		
38		Sonic	100	1175.0	56-60" Wet, dark gray, medium to coarse SAND, strong hydrocarbon odor		2-inch ID Well Screen (28-48')
39							0.02-inch slot PVC
40			 	209.0	0-60" Wet, brown medium to fine SAND, some orange Silt lenses, some black staining, slight	※ ※	
42					hydrocarbon odor		
43				19.5	60-100" Wet, orange/brown medium to coarse SAND, trace fine round Gravel	∰ ∭	
45		Sonic	83				
46 47				<u> </u>			
48	MW-7 (48-50)					1000 A	
49	111 11 -7 (40-30)		1		N. P	┨ └┴──┼	2' Sump (48-50')
50 51		Sonic	0	NA	No Recovery		PVC End Cap (50')
52						1	
53	SAMPLING METH	HOD			End of boring at 53' bgs NAPL observed 35' - 40' bgs		
	SS = SPLIT SPOON				Hard drilling from 10-20 ft bgs. Casing advanced to 20 ft bgs. Drilled through rock from 50 to 53 ft bgs. No rock re	trieved from 50 to 53 ft bgs	
	A = AUGER CUTTI C = CORED	NGS			Hand cleared to 5' bgs, sonic drilled from 5' to 53' bgs		
	WH = WEIGHT OF	HAMMER (RODS)					

					PARSONS	BORING/WELL ID	: MW-8
					DRILLING RECORD		Sheet 1 of 1
Contracto	or: Aquifer Dri	lling & Testing				Location Description:	
Driller:	Greg Rivera	ı			PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	On MTA/Metro North Pro	
Inspector	: Matt Bruno				PROJECT NUMBER: 446110-04000	40' west of railroad track, Sugar Refining southern p	
Rig Type:		nted Sonic Rig					
		WATER OBSERVAT	ΓIONS			Location Plan	
Water	DTW	DTW			Weather: Partly cloudy, 40s		
Level	~6' bgs	4.68					
Date	10-18-2012	11-9-12			Date/Time Start: 11-06-2012/1345	See Site Plan	
Time	1430	1322					
Meas.	ft bgs -	Top of Casing			Date/Time Finish: 11-06-2012/1400		
From	Soil cuttings	o.m.m.		nvn		CONTRACTOR O	001111mm
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.		(%)	(ppm)			
+2							Y 11 Y 1
+1 0		Vac-tron	NA	0.0	0-12" Dry, gray fine to medium SAND and sub-angular to sub-round GRAVEL		Locking J-plug on casing Flushmount manhole
1		Vac-tron	NA NA	0.0	12"-5.5' Dry, dark brown, fine to coarse SAND, some Cobble, little Gravel, little Boulder, trace Brick		Fiushmount mannoie
2		Vac-tron	NA NA	0.0	and Silt		Grout (0-2')
3		Vac-tron	NA	0.0	5.5'-6' Moist, dark brown, fine to medium SAND and BOULDER, little Silt	1 (3) (3)	Glour (0-2)
4		Vac-tron	NA	0.0	,		2" PVC riser (0.5-4')
5		vac-tron	IVA	0.0	0-24" Moist, tan/brown SILT, some sub-angular Gravel, little round Gravel	3 3 -	2 1 ve liser (0.3-4)
6		1		3.7	24-36" Moist-wet, black/dark gray sub-angular to sub-round COBBLE/GRAVEL, some Silt and fine to		
7		Sonic	60		coarse Sand, hydrocarbon odor		
8		1				8 - 8	
9	MW-8 (9.5-10)						2" ID PVC well screen
10				3.8	0-12" Wet, soft, black/dark brown SILT, little fine to medium Sand, trace sub-round Gravel		0.002" slot (4-14')
11				0.0	12-30" Wet, dark brown/black fine to medium SAND, some Silt, little sub-angular Cobble/Gravel	38 88	
12				0.0	30-66" Moist, soft, red/brown fine to medium SAND, little Silt, trace Clay		#2 silica (2-14')
13							
14		Sonic	46				
15		Joine	40				PVC end cap 14'
16		1					
17		1					
18		4					
19			_		La activity and the state of th	4	
20		4			0-20" Wet, soft. brown/red fine to medium SAND, little Silt		
21		Comin	60	0.0	20-36" Dry-moist, brown/red SILT, little fine Sand, orange/tan Silt lenses, trace sub-round Gravel		
22		Sonic	60	0.0			
23	MW-8 (24.5-25	1					
24 25	141 44-0 (24.3-23	1		1	End of bosing of 25' box	-	
25	SAMPLING MET	HOD			End of boring at 25' bgs Hand cleared to 5' bgs, sonic drilled from 5' to 25' bgs.		
	SS = SPLIT SPOON				Trans course of 5 ogs, some timed from 5 to 25 ogs.		
	A = AUGER CUTT						
	C = CORED	1105					
		HAMMER (RODS)					

					PARSONS	BORING/WELL II	D: MW-9
_					DRILLING RECORD		Sheet 1 of 1
Contracto	or: Aquifer Dril	ling & Testing				Location Description:	
Driller:	Greg Rivera				PROJECT NAME: Consolidated Edison - Former Ludlow Street Works	Located approximately 30	
Inspector	: Matt Bruno				PROJECT NUMBER: 446110-04000	and approximately 50' so Refining southern proper	
Rig Type:	Track Moun	ited Sonic Rig			-		
8 71		VATER OBSERVATI	ONS			Location Plan	
Water	DTW	DTW			Weather: Clear, 40s		
Level		3.89					
Date		11-9-12			Date/Time Start: 11-07-2012/0845	See Site Plan	
Time		1135			D. (70° - 10° - 11 07 2012/1005		
Meas. From		Top of Casing			Date/Time Finish: 11-07-2012/1005	4	
Sample	Location/	SPT	Rec.	PID	FIELD IDENTIFICATION OF MATERIAL	SCHEMATIC	COMMENTS
Depth	Sample I.D.		(%)	(ppm)			
+2							
+1						ļ <u></u>	Locking J-plug on casing
0		Vac-tron/Sonic	onic NA 12"-5' Dry, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 0.0 round Gravel onic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-anguonic NA 5-5.5' Moist, brown/black fine to coarse sub-anguonic NA 5-5.5' Moist, b		0-12": Dry, brown/gray fine to medium SAND and sub-angular to sub-round GRAVEL, trace Metal	W # #	Flushmount manhole
1		Vac-tron/Sonic		12"-5' Dry, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-angular to round Gravel 5-5.5' Moist, brown/black fine to coarse SAND and COBBLE, some fine to coarse sub-angular to round Gravel		1 20 2	C+ (0.2h
3		Vac-tron/Sonic Vac-tron/Sonic		0.0			Grout (0-2')
4		Vac-tron/Sonic	NA NA	 			2" PVC riser (0.5-4')
5		. ac tron/bonic	1121		No Recovery	1 3 3 1	2 1 (0.34)
6	1	1					
7		Sonic	0	NA			
8							
9	MW-9 (9-9.5)						2" ID PVC well screen
10					0-12" Wet, dark brown/black GRAVEL, some sub-round Boulders and Cobbles, little fine Sand and	98 83	0.002" slot PVC (4-14')
11		G	20	2.7	Silt	1 8848	
12		Sonic	20	3.7			#2 silica (2-14')
13 14							
15				0.5	0-12" Wet, dark brown/black GRAVEL, some sub-round Boulders and Cobbles, little fine Sand and	⊗⊨⊗	PVC end cap 14'
16		1		0.5	Silt, trace Metal wire fragment		1 ve cha cap 1 v
17		Sonic	60	1.6	12-36" Moist, black SILT and fine SAND, little Shell fragments		
18							
19							
20					0-36" Wet, very soft, red/brown fine SAND, some Silt		
21			60	0.0			
22		Sonic	60	0.0			
24							
25			<u> </u>	0.0	0-12" Wet, very soft, red/brown SILT, some fine Sand	1	
26		1		0.3	12-24" Moist, medium dense, red/brown SILT, some Clay		
27		Sonic	90	0.0	24-36" Moist, red/brown SILT and sub-round GRAVEL		
28]		0.0	36-54" Wet, red/tan, medium to coarse SAND, little sub-round Gravel, little Silt		
29			ļ			4	
30				0.0	0-30" Wet, red/tan, medium to coarse SAND, some sub-angular Cobble and Boulder, little sub-round		
31		Sonic	66	0.2	Gravel, little Silt 30-35" Moist, red/orange SILT, trace sub-round Gravel		
33		Some	00	0.2	35-40" Wet, red/orange medium SAND, little Silt, trace sub-round Cobble/Gravel		
34		1		0.0			
35	MW-9(35-35.5)		<u> </u>	<u> </u>	0-12" Wet, red/orange medium SAND, little Silt, trace sub-round Cobble/Gravel	1	
36	1	1			12-24" Moist, med-dense, red/orange fine SAND, some Silt, trace sub-round Gravel		
37		Sonic	80	0.0	24-30" Moist, med-dense, red/orange fine SAND, some Silt, trace sub-angular Gravel		
38					30-48" Dry, red/brown SILT, little white medium Sand lenses		
39]		4	
40	CAMPA PAGES	HOD			End of boring at 40' bgs	<u> </u>	<u> </u>
	SAMPLING METI				Hand cleared to 5' bgs, sonic drilled from 5' to 40' bgs		
	SS = SPLIT SPOON A = AUGER CUTTI						
	C = CORED						
		HAMMER (RODS)					

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 11-26-12 Sampling Date: 11-26-12 Samplers: Zohar Lavy Parsons / Somerset, NJ SAMPLE ID: MW-1 Sampling Method: Low flow purge (Monsoon Pump) WELL PURGING Static Water Level (TOC): 14.12 Depth to Well Bottom (TOC): 20.20 **CALCULATIONS:** Ft. of Water in Well X (GAL / FT) =Gallons 2-inch Casing: Ft. of Water in Well Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 = Gallons 4-inch Casing: Gallons Ft. of Water in Well x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: No Odor Other: Clear **FIELD TESTS PURGE PURGE PURGE PURGE PURGE PURGE PURGE PURGE** Time 0940 0950 0955 1005 1010 1015 1020 1025 Depth To Water (TOC) (ft) 18.90 Below Pump Depth To Pump (TOC) (ft) 19.20 19.20 19.20 19.20 19.20 19.20 19.20 19.20 Flow Rate (ml/min) ~100 ~100 ~100 ~100 ~100 ~250 ~100 ~100 Volume of Water Purged ~0.75 ~1.75 ~2.25 ~0.5 ~1 ~1.25 ~1.5 ~2 7.03 7.00 7.01 7.01 7.02 7.03 7.03 7.03 pH (s.u.) Conductivity (mS/cm) 2.68 2.63 2.57 2.54 2.49 2.36 2.50 2.48 Turbidity (NTUs) 495 740 548 226 118 54 36.7 12.9 9.48 9.44 9.79 Dissolved Oxygen (mg/L) 13.37 9.61 9.70 10.03 10.03 Temperature (Degrees C) 16.49 16.67 15.66 15.86 16.00 16.03 16.06 16.35 ORP (mV) 89 87 100 46 33 19 18 Salinity (%) 0.9 0.9 0.9 0.9 0.9 0.9 0.9 0.9 TDS (g/L) 1.73 1.71 1.68 1.64 1.62 1.60 1.60 1.59 **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory

Sample collected at 1030; purged approximately 2.5 gallons

Other Notes:

GROUNDWATER SAMPLING RECORD

SITE NAME:	Con Edison (Ludlow)			
PROJECT NUMBER:	446110-04000			
Purge Date:	11-26-12			
Sampling Date:	11-26-12		-	
Samplers:	Zohar Lavy	of	Parsons / Somerset, NJ	
Campioro			Taloution Completel, 140	
SAMPLE ID:	MW-1			
Sampling Method:	Low flow purge (Monsoon Pu	mp)		
campanig memera		···F/		
WELL PURGING				
Static Water Level (TOC):	14.12			
Depth to Well Bottom (TOC):	20.20			
CALCULATIONS:	Ft. of Water in Well	X (GAL / FT) =	Gallons	
2-inch Casing:	Ft. of Water in Well	x 0.16 =	Gallons	
3-inch Casing:	Ft. of Water in Well	x 0.32 =	Gallons	
4-inch Casing:	Ft. of Water in Well	x 0.64 =	Gallons	
Method:	Low Flow Pump			
SAMPLE DESCRIPTION				
Odor:	No Odor			
Other:	Clear			
FIELD TESTS				
	SAMPLE			
Time	1030			
Depth To Water (TOC) (ft)	Below Pump			
Depth To Pump (TOC) (ft)	19.20			
Flow Rate (ml/min)	~100			
Volume of Water Purged	~2.50			
pH (s.u.)	7.03			
Conductivity (mS/cm)	2.43			
Turbidity (NTUs)	8.6			
Dissolved Oxygen (mg/L)	10.05			
Temperature (Degrees C)	16.34			
ORP (mV)	17			
Salinity (%)	0.9			
TDS (g/L)	1.59			
SAMPLE ANALYSIS / LABORATO	DRY			
Analyze For:	TCL VOC's , TCL SVOCs, TA	AL Metals, Cyanide		
Shipped Via:	Chemtech			
Laboratory				
Other Notes:	Sample collected at 1030; pu	rged approximately 2.5 gallons	· · · · · · · · · · · · · · · · · · ·	
	·			

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 11-27-12 Sampling Date: 11-27-12 Samplers: Zohar Lavy Parsons / Somerset, NJ SAMPLE ID: MW-3 Sampling Method: Low flow purge (Monsoon Pump) WELL PURGING Static Water Level (TOC): 16.58 Depth to Well Bottom (TOC): 29.90 **CALCULATIONS:** Ft. of Water in Well X (GAL / FT) = Gallons 2-inch Casing: Ft. of Water in Well Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 = Gallons 4-inch Casing: Gallons Ft. of Water in Well x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: Strong hydrocarbon odor Other: Clear, Sheen **FIELD TESTS PURGE PURGE PURGE PURGE PURGE PURGE PURGE PURGE** Time 1430 1435 1440 1445 1450 1455 1500 1505 Depth To Water (TOC) (ft) 17.25 16.82 16.80 16.78 16.79 16.78 16.79 16.81 Depth To Pump (TOC) (ft) 26.00 26.00 26.00 26.00 26.00 26.00 26.00 26.00 Flow Rate (ml/min) ~300 ~350 ~350 ~300 ~350 ~300 ~350 ~350 Volume of Water Purged ~0.5 ~1.0 ~1.5 ~2.25 ~3.0 ~3.5 ~4.25 ~5.0 7.22 7.21 7.21 7.20 7.20 7.20 7.20 7.19 pH (s.u.) Conductivity (mS/cm) 23.7 15.25 27.3 20.2 18.5 14.65 13.31 13.18 Turbidity (NTUs) 572 393 114 17.9 10.5 5.7 4.2 3.1 Dissolved Oxygen (mg/L) 14.70 13.25 12.36 11.63 10.98 10.11 9.86 9.79 Temperature (Degrees C) 13.16 11.23 10.12 10.61 11.90 12.03 12.16 13.80 ORP (mV) -57 -53 -50 -49 -49 -49 -50 -50 Salinity (%) 13.8 11.2 10.5 9.4 8.1 6.9 6.5 6.0 TDS (g/L) 15.7 14.1 12.7 11.6 8.19 8.07 8.0 **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1510; purged approximately 5.5 gallons; NAPL in purge water

GROUNDWATER SAMPLING RECORD

PROJECT NUMBER:				
	446110-04000			
Purge Date:	11-27-12			
Sampling Date:	11-27-12			
Samplers:	Zohar Lavy	of	Parsons / Somerset, NJ	
SAMPLE ID:	MW-3			
Sampling Method:	Low flow purge (Monsoon Pu	ımp)		
WELL PURGING				
Static Water Level (TOC):	16.58			
Depth to Well Bottom (TOC):	29.90			
CALCULATIONS:	Ft. of Water in Well	X (GAL / FT) =	Gallons	
2-inch Casing:	Ft. of Water in Well	x 0.16 =	Gallons	
3-inch Casing:	Ft. of Water in Well	x 0.32 =	Gallons	
4-inch Casing:	Ft. of Water in Well	x 0.64 =	Gallons	
Method:	Low Flow Pump			
SAMPLE DESCRIPTION				
Odor:	Strong hydrocarbon odor			
Other:	Clear, Sheen			
FIELD TESTS	0.00., 0.100.1			
	SAMPLE			
Time	1510			
Depth To Water (TOC) (ft)	16.81			
Depth To Pump (TOC) (ft)	26.00			
Flow Rate (ml/min)	~350			
Volume of Water Purged	~5.5			
pH (s.u.)	7.19			
Conductivity (mS/cm)	13.00			
Turbidity (NTUs)	1.1			
Dissolved Oxygen (mg/L)	9.32			
Temperature (Degrees C)	14.18			
ORP (mV)	-50			
Salinity (%)	6.0			
TDS (g/L)	7.91			
100 (g/L)	7.51			
SAMPLE ANALYSIS / LABORATO				
Analyze For:	TCL VOC's , TCL SVOCs, TA	AL Metals, Cyanide		
Shipped Via:	Chemtech			
Laboratory				
Other Notes:	0	rged approximately 5.5 gallons	NADI is sures water	

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 11-27-12 Sampling Date: 11-27-12 Samplers: Parsons / Somerset, NJ Zohar Lavy SAMPLE ID: MW-4 Sampling Method: Low flow purge (Monsoon Pump) WELL PURGING Static Water Level (TOC): 14.30 Depth to Well Bottom (TOC): 18.35 **CALCULATIONS:** Ft. of Water in Well X (GAL / FT) =2-inch Casing: Ft. of Water in Well Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 = Gallons 4-inch Casing: Gallons Ft. of Water in Well x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: No odor Other: Very turbid **FIELD TESTS PURGE PURGE PURGE PURGE PURGE** SAMPLE Time 1120 1125 1130 1140 1145 1150 Depth To Water (TOC) (ft) 16.22 Below Pump Dry Below Pump Dry Below Pump Depth To Pump (TOC) (ft) 17.35 17.35 17.35 17.35 17.35 17.35 Flow Rate (ml/min) ~250 ~350 NA ~100 ~100 NA Volume of Water Purged ~0.25 ~0.75 NA ~1.0 NA ~1.25 6.42 7.37 NA 7.85 NA 7.56 pH (s.u.) Conductivity (mS/cm) 2.07 NA NA 2.61 1.87 1.62 Turbidity (NTUs) Error Error NA Error NA Error 24.73 Dissolved Oxygen (mg/L) 23.4 NA 24.33 NA 24.11 Temperature (Degrees C) 12.13 11.88 NA 10.67 NA 11.36 ORP (mV) 163 158 NA 170 NA 152 Salinity (%) 8.0 0.6 NA 0.6 NA 0.6 TDS (g/L) 1.62 1.32 NA 1.18 **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1150; purged approximately 1.25 gallons; well ran dry two times prior to sampling

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 11-27-12 Sampling Date: 11-27-12 Samplers: Zohar Lavy Parsons / Somerset, NJ SAMPLE ID: MW-7 Sampling Method: Low flow purge (Monsoon Pump) WELL PURGING Static Water Level (TOC): 39.32 Depth to Well Bottom (TOC): 51.77 **CALCULATIONS:** Ft. of Water in Well X (GAL / FT) = Gallons 2-inch Casing: Ft. of Water in Well Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 = Gallons 4-inch Casing: Gallons Ft. of Water in Well x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: Slight hydrocarbon odor Other: Clear, Sheen **FIELD TESTS PURGE PURGE PURGE PURGE PURGE PURGE PURGE** SAMPLE Time 1240 1245 1250 1255 1300 1305 1310 1315 Depth To Water (TOC) (ft) 39.45 39.45 39.46 39.47 39.45 39.46 39.45 39.45 Depth To Pump (TOC) (ft) 48.77 48.77 48.77 48.77 48.77 48.77 48.77 48.77 Flow Rate (ml/min) ~350 ~300 ~300 ~300 ~250 ~350 ~300 ~300 Volume of Water Purged ~0.5 ~1.0 ~1.5 ~2.0 ~2.5 ~3.0 ~3.5 ~4.0 7.21 7.18 7.17 7.16 7.16 7.17 7.17 7.18 pH (s.u.) Conductivity (mS/cm) 3.84 3.84 3.82 3.81 3.85 3.81 3.80 3.80 Turbidity (NTUs) 482 303 212 165 92.5 33.2 20.6 17.5 Dissolved Oxygen (mg/L) 11.73 10.27 9.31 8.08 7.34 7.17 7.04 6.98 Temperature (Degrees C) 14.49 14.61 14.52 14.55 14.63 14.59 14.47 14.51 ORP (mV) 0 -1 -3 -4 -4 -4 -4 Salinity (%) 1.4 1.4 1.4 1.4 1.4 1.4 1.4 TDS (g/L) 2.46 2.45 2.45 2.44 2.44 2.43 2.43 2.42 **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1315; purged approximately 4.0 gallons

GROUNDWATER SAMPLING RECORD

SITE NAME:	Con Edison (Ludlow	')						
PROJECT NUMBER:	446110-04000							
Purge Date:	11-26-12							
Sampling Date:	11-26-12							
Samplers:	Zohar L	.avy	of		Parsor	s / Somerset	, NJ	
SAMPLE ID:	MW-8 and MW-18 (c	duplicate)						
Sampling Method:	Low flow purge (Mor							
Camping memori		iocon i amp)						
WELL PURGING								
Static Water Level (TOC):	5.27							
Depth to Well Bottom (TOC):	13.15							
CALCULATIONS:	Ft. of Water in Well		X (GAL / FT) =		Gallons			
2-inch Casing:	Ft. of Water in Well		x 0.16 =		Gallons			
3-inch Casing:	Ft. of Water in Well		x 0.32 =		Gallons			
4-inch Casing:	Ft. of Water in Well		x 0.64 =		Gallons			
Method:	Low Flow Pump							
SAMPLE DESCRIPTION								
Odor:	No Odor							
Other:	Clear							
FIELD TESTS	PURGE	PURGE	PURGE	PURGE	PURGE	PURGE	PURGE	PURGE
Time	1205	1215	1220	1225	1230	1235	1240	1245
Depth To Water (TOC) (ft)	5.38	5.45	5.45	5.45	5.45	5.46	5.46	5.47
Depth To Water (TOC) (ft) Depth To Pump (TOC) (ft)	12.15	12.15	12.15	12.15	12.15	12.15	12.15	12.15
Flow Rate (ml/min)	~150	~150	~250	~200	~250	~250	~250	~250
Volume of Water Purged	~0.25	~0.5	~1.0	~1.25	~1.75	~2.25	~2.5	~2.75
pH (s.u.)	7.05	7.00	6.96	6.98	6.98	6.98	6.99	6.99
Conductivity (mS/cm)	4.84	4.70	4.60	4.59	4.59	4.59	4.59	4.59
Turbidity (NTUs)	Error	Error	Error	769	503	411	282	146
Dissolved Oxygen (mg/L)	11.69	10.54	9.15	8.82	8.73	8.65	8.48	8.04
Temperature (Degrees C)	14.59	16.04	17.05	17.06	17.03	17.14	17.29	17.46
ORP (mV)	176	147	119	110	107	102	95	89
Salinity (%)	1.8	1.8	1.7	1.7	1.7	1.7	1.7	1.7
TDS (g/L)	3.11	2.99	2.94	2.94	2.94	2.93	2.94	2.94
= (3, _)	3							
SAMPLE ANALYSIS / LABORATO	ORY							
Analyze For:	TCL VOC's, TCL S\	/OCs, TAL Met	als, Cyanide					
•			•					
	<u> </u>							
Shipped Via:	Chemtech							
Laboratory								
Other Notes:	Sample collected at	4000						

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 11-26-12 Sampling Date: 11-26-12 Samplers: Zohar Lavy Parsons / Somerset, NJ SAMPLE ID: MW-8 and MW-18 (duplicate) Sampling Method: Low flow purge (Monsoon Pump) WELL PURGING Static Water Level (TOC): 5.27 Depth to Well Bottom (TOC): 13.15 **CALCULATIONS:** Ft. of Water in Well X (GAL / FT) =Gallons 2-inch Casing: Ft. of Water in Well Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 = Gallons 4-inch Casing: Gallons Ft. of Water in Well x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: No Odor Other: Clear **FIELD TESTS PURGE PURGE PURGE PURGE PURGE PURGE** SAMPLE Time 1250 1255 1300 1305 1310 1315 1320 Depth To Water (TOC) (ft) 5.47 5.47 5.7 5.48 5.47 5.47 5.47 Depth To Pump (TOC) (ft) 12.15 12.15 12.15 12.15 12.15 12.15 12.15 Flow Rate (ml/min) ~250 ~250 ~350 ~200 ~300 ~300 ~250 Volume of Water Purged ~4.5 ~3.25 ~3.75 ~4.75 ~5.25 ~5.75 ~6.25 6.99 6.99 6.99 6.99 6.99 6.99 6.99 pH (s.u.) Conductivity (mS/cm) 4.58 4.62 4.64 4.60 4.63 4.61 4.58 35.8 Turbidity (NTUs) 137 161 98.9 73.2 59.4 43.6 7.28 Dissolved Oxygen (mg/L) 7.83 7.51 7.31 7.19 7.12 7.01 17.87 Temperature (Degrees C) 17.60 17.83 17.86 17.88 17.89 17.91 ORP (mV) 86 85 84 85 84 83 82 Salinity (%) 1.7 1.7 1.7 1.7 1.7 1.7 1.7 TDS (g/L) 2.95 2.94 2.95 2.94 2.94 2.94 2.94 **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1320; purged approximately 6.25 gallons

GROUNDWATER SAMPLING RECORD

SITE NAME:	Con Edison (Ludlow	/)						
PROJECT NUMBER:	446110-04000							
Purge Date:	11-26-12							
Sampling Date:	11-26-12							
Samplers:	Zohar L	.avy	of		Parsor	s / Somerset	, NJ	
SAMPLE ID:	MW-9, MW-9MS, M\	W-9MSD						
Sampling Method:	Low flow purge (Mor							
, 3		1,						
WELL PURGING								
Static Water Level (TOC):	4.45							
Depth to Well Bottom (TOC):	12.15							
CALCULATIONS:	Ft. of Water in Well		X (GAL / FT) =		Gallons			
2-inch Casing:	Ft. of Water in Well		x 0.16 =		Gallons			
3-inch Casing:	Ft. of Water in Well		x 0.32 =	<u></u>	Gallons			
4-inch Casing:	Ft. of Water in Well		x 0.64 =		Gallons			
Method:	Low Flow Pump							
SAMPLE DESCRIPTION								
Odor :	Slight hydrocarbon o	dor						
Other:	Clear							
FIELD TESTS	5,15,05	55.05	5115.05		5.15.05	5115.05	5115.05	0
_	PURGE	PURGE	PURGE	PURGE	PURGE	PURGE		SAMPLE
Time	1410	1415	1420	1425	1430	1435	1440	1445
Depth To Water (TOC) (ft)	4.49	4.49	4.49	4.52	4.51	4.51	4.51	4.52
Depth To Pump (TOC) (ft)	11.15	11.15	11.15	11.15	11.15	11.15	11.15	11.15
Flow Rate (ml/min)	~350	~350	~350	~350	~350	~300	~300	~350
Volume of Water Purged	~1.0	~1.5	~2.0	~2.5	~3.0	~3.5	~4.0	~4.5
pH (s.u.)	7.29	7.29	7.29	7.28	7.29	7.27	7.27	7.26
Conductivity (mS/cm)	9.59	9.62	9.74	9.83	9.90	9.88	9.83	9.81
Turbidity (NTUs)	32.7	8.8	0	0	0	0	0	0
Dissolved Oxygen (mg/L)	9.56	7.45	6.6	6.24	6.01	5.89	5.81	5.76
Temperature (Degrees C)	15.07	14.92	14.89	14.90	14.88	14.72	14.71	14.74
ORP (mV)	-133	-133	-133	-133	-134	-134	-133	-134
Salinity (%)	4.2	4.3	4.3	4.4	4.3	4.4	4.3	4.3
TDS (g/L)	6.04	6.08	6.15	6.19	6.23	6.22	6.21	6.21
SAMPLE ANALYSIS / LABORATO	D V							
		OCa TAL Mar	ala Cuanida					
Analyze For:	TCL VOC's, TCL S\	/UUS, TAL Met	ais, Cyanide					
	-							
Shipped Via:	Chemtech							
Laboratory	Onomidon							
Laboratory								
Other Notes:	Sample collected at	1.445: purgod a	poprovimatoly 4 F	gallons				
Other Notes:	Sample collected at	1740, purged a	ippioximalely 4.5	yalluris				

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 4-5-13 Sampling Date: 4-8-13 Samplers: Zohar Lavy Parsons / Somerset, NJ SAMPLE ID: MW-1 Sampling Method: Low flow purge (Monsoon Pump) **WELL PURGING** Static Water Level (TOC): 13.36 Depth to Well Bottom (TOC): 20.30 **CALCULATIONS:** Ft. of Water in Well X (GAL/FT) =Gallons 2-inch Casing: Ft. of Water in Wel Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 =Gallons 4-inch Casing: Ft. of Water in Well Gallons x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: No odor Other: Very turbid **FIELD TESTS PURGE PURGE PURGE PURGE PURGE PURGE** SAMPLE Time 1215 1225 1245 1300 1210 1220 1240 Depth To Water (TOC) (ft) 16.25 Below Pump 17.8 17.22 17.79 17.4 Dry Depth To Pump (TOC) (ft) 19.3 19.3 19.3 19.3 19.3 19.3 19.3 ~150 ~100 ~100 ~100 ~100 Flow Rate (ml/min) NA NA Volume of Water Purged ~0.75 ~1.75 ~0.5 ~1 ~1 ~1.5 ~1.5 7.09 7.17 7.09 NA 7.23 NA 7.18 pH (s.u.) Conductivity (mS/cm) 1.98 1.98 NA 1.89 NA 1.88 1.96 Turbidity (NTUs) 166 122 125 NA 130 NA 122 Dissolved Oxygen (mg/L) 9.15 3.36 3.27 NA 5.03 NA 3.44 14.97 Temperature (Degrees C) 14.46 15.22 NA 15.6 NA 15.84 77 88 89 89 45 NA 41 ORP (mV) 1.0 1.0 1.0 NA 1.0 NA 1.0 Salinity (%) NA 1.27 1.27 1.21 1.20 TDS (g/L) 1.26 NA **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide, dissolved Metals Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1300; purged approximately 1.75 gallons; well ran dry several times prior to sampling

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 4-8-13 Sampling Date: 4-8-13 Samplers: Zohar Lavy, Dennis Miller Parsons / Somerset, NJ SAMPLE ID: Sampling Method: Low flow purge (Monsoon Pump) **WELL PURGING** Static Water Level (TOC): 10.95 Depth to Well Bottom (TOC): 19.30 **CALCULATIONS:** Ft. of Water in Well X (GAL/FT) =Gallons 2-inch Casing: Ft. of Water in Wel Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 =Gallons 4-inch Casing: Ft. of Water in Well Gallons x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION No Odor Odor: Other: Clear **FIELD TESTS PURGE PURGE PURGE** PURGE **PURGE PURGE** SAMPLE Time 0752 0757 0802 0807 0812 0817 0822 Depth To Water (TOC) (ft) 11.50 12.30 11.87 12.15 12.40 12.45 12.35 Depth To Pump (TOC) (ft) 15.0 15.0 15.0 15.0 15.0 15.0 15.0 100 250 250 250 Flow Rate (ml/min) 100 200 250 Volume of Water Purged 0.125 0.25 0.375 0.75 1.25 0.5 1.0 5.92 6.04 5.61 5.79 5.96 6.0 6.06 pH (s.u.) Conductivity (mS/cm) Error Error Error Error Error Error Error Turbidity (NTUs) 162 34.2 12.0 3.4 0.0 0.0 0.0 Dissolved Oxygen (mg/L) 2.96 2.45 2.06 1.66 1.44 1.34 1.24 10.64 10.78 Temperature (Degrees C) 10.02 10.05 10.10 10.39 10.62 110 60 40 23 8 -2 -11 ORP (mV) Error Error Error Error Error Salinity (%) Error Error TDS (g/L) Error Error Error Error Error Error Error SAMPLE ANALYSIS / LABORATORY Analyze For: TCL VOC's, TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 0822; purged approximately 1.25 gallons

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow)

PROJECT NUMBER: 446110-04000

Purge Date: 4-8-13

Sampling Date: 4-8-13

Samplers: Zohar Lavy, Dennis Miller of Parsons / Somerset, NJ

SAMPLE ID: MW-3 and duplicate

Sampling Method: Low flow purge (Monsoon Pump)

WELL PURGING

 Static Water Level (TOC):
 16.40

 Depth to Well Bottom (TOC):
 29.95

CALCULATIONS: Ft. of Water in Well X (GAL/FT) =Gallons 2-inch Casing: Ft. of Water in Wel Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 =Gallons 4-inch Casing: Ft. of Water in Well Gallons x 0.64 =

Method: Low Flow Pump

SAMPLE DESCRIPTION

Odor: Strong hydrocarbon odor

Other: Clear, Sheen, NAPL Blebs

FIELD TESTS

PURGE PURGE PURGE PURGE **PURGE PURGE** SAMPLE Time 0924 0929 0934 0939 0944 0949 0954 Depth To Water (TOC) (ft) 17.05 16.95 16.90 16.91 16.92 16.89 16.90 Depth To Pump (TOC) (ft) 25.00 25.00 25.00 25.00 25.00 25.00 25.00 Flow Rate (ml/min) 250 250 250 250 250 250 250 Volume of Water Purged 0.25 0.5 1.0 1.25 1.5 1.75 2.0 6.95 6.96 6.97 6.96 pH (s.u.) 6.73 6.9 6.92 Conductivity (mS/cm) 37.1 86 46.9 33.7 31.6 28.6 30.7 Turbidity (NTUs) 105 46.8 38.3 37.9 30.0 32.2 31.1 Dissolved Oxygen (mg/L) 7.94 8.24 7.37 6.66 5.98 5.60 5.44 Temperature (Degrees C) 14.20 14.90 14.96 15.01 15.02 14.72 14.49 -26 -26 -31 -30 -31 -28 -26 ORP (mV) 57.2 29.1 22.9 20.8 18.1 17.6 16.9 Salinity (%)

27.5

SAMPLE ANALYSIS / LABORATORY

TDS (g/L)

Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide

49.1

22.3

18.3

20.5

17.1

17.8

Shipped Via: Chemtech
Laboratory

Other Notes: Sample collected at 0954; purged approximately 2 gallons; NAPL in purge water

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow)

PROJECT NUMBER: 446110-04000

 Purge Date:
 4-8-13

 Sampling Date:
 4-8-13

Samplers: Zohar Lavy, Dennis Miller of Parsons / Somerset, NJ

SAMPLE ID: MW-4

Sampling Method: Low flow purge (Monsoon Pump)

WELL PURGING

Static Water Level (TOC): 14.30

Depth to Well Bottom (TOC): 18.35

CALCULATIONS: Ft. of Water in Well X (GAL/FT) =Gallons 2-inch Casing: Ft. of Water in Wel Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 =Gallons 4-inch Casing: Ft. of Water in Well Gallons x 0.64 =

Method: Low Flow Pump

SAMPLE DESCRIPTION

Odor: No odor

Other: Very turbid

FIELD TESTS

PURGE PURGE PURGE PURGE PURGE SAMPLE Time 0750 0755 0800 0810 0815 0840 Below Pump Depth To Water (TOC) (ft) 16.4 Dry Dry Dry Dry Depth To Pump (TOC) (ft) 17.4 17.4 17.4 17.4 17.4 17.4 ~100 ~100 ~100 ~100 ~100 ~100 Flow Rate (ml/min) Volume of Water Purged ~0.25 NA NA NA NA ~0.75 8.12 NA NA NA NA 8.05 pH (s.u.) Conductivity (mS/cm) 4.91 NA NA NA NA 4.86 Turbidity (NTUs) Error NA NA NA NA Error Dissolved Oxygen (mg/L) 13.58 NA NA NA NA 9.87 Temperature (Degrees C) 11.06 NA NA NA NA 10.67 4.48 NA NA NA NA 169 ORP (mV) 2.6 NA NA 2.6 NA NΑ Salinity (%)

NA

SAMPLE ANALYSIS / LABORATORY

TDS (g/L)

Analyze For: TCL VOC's , TCL SVOCs, TAL Metals, Cyanide, dissolved Metals

3.14

NA

NA

3.12

Shipped Via: Chemtech
Laboratory

NA

Other Notes: Sample collected at 0840; purged approximately 0.75 gallons; well ran dry several times prior to sampling

GROUNDWATER SAMPLING RECORD

	Con Edison (Ludi	OW)						
PROJECT NUMBER:	446110-04000							
Purge Date:	4-5-13		<u>—</u>					
Sampling Date:	4-5-13							
Samplers:	Zohai	r Lavy	of		Parson	s / Somerset	, NJ	
SAMPLE ID:	MW-5, MW-5MS,	MW-5 MSD						
Sampling Method:	Low flow purge (N	fonsoon Pump)						
WELL PURGING								
Static Water Level (TOC):	12.05							
Depth to Well Bottom (TOC):	21.30	_"						
CALCULATIONS:	Ft. of Water in We	_ ell	X (GAL / FT) =	=	Gallons			
2-inch Casing:	Ft. of Water in We	el	x 0.16 =		Gallons			
3-inch Casing:	Ft. of Water in We	ell	x 0.32 =		Gallons			
4-inch Casing:	Ft. of Water in We	ell	x 0.64 =		Gallons			
Method:	Low Flow Pump				_			
SAMPLE DESCRIPTION								
Odor:	Cliabt budrooorbo	n Odor						
Odor:	Slight hydrocarbo	n Odor						
FIELD TESTS	Clear							
FIELD IESIS	PURGE	DUDGE	DUDGE	DUDGE	DUDGE	DUDGE	DUDGE	CAMPLE
Time		PURGE	PURGE	PURGE	PURGE	PURGE	1405	SAMPLE
	1335	1340	1345	1350	1355	1400		1410
Depth To Water (TOC) (ft)	12.7	12.73	12.72	12.72	12.75	12.73	12.72	12.73
Depth To Pump (TOC) (ft)	20.3	20.3	20.3	20.3	20.3	20.3	20.3	20.3
Flow Rate (ml/min)	~300	~350	~350	~350	~350	~350	~350	~350
Volume of Water Purged	~0.75	~1.5	~2.25	~3.0	~3.75	~3.5	~5.0	~5.5
pH (s.u.)	6.91	6.94	6.98	6.98	6.98	6.99	6.99	6.99
Conductivity (mS/cm)	Error	Error	Error	Error	Error	Error	Error	Error
Turbidity (NTUs)	117	165	125	102	90.1	62.5	51.3	32.6
Dissolved Oxygen (mg/L)	5.29	1.53	0.99	0.87	0.82	0.77	0.75	0.73
Temperature (Degrees C)	13.93	13.65	13.71	13.72	13.75	13.78	13.74	13.76
ORP (mV)	-40	-49	-60	-61	-62	-64	-65	-66
Salinity (%)	Error	Error	Error	Error	Error	Error	Error	Error
TDS (g/L)	Error	Error	Error	Error	Error	Error	Error	Error
SAMPLE ANALYSIS / LABORATO	DRY							
Analyze For:	TCL VOC's, TCL	SVOCs, TAL Me	etals, Cyanide					
Shipped Via:	Chemtech							
Laboratory	Onomicon							
Laboratory								
·								

GROUNDWATER SAMPLING RECORD

SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 4-8-13 Sampling Date: 4-8-13 Samplers: Zohar Lavy, Dennis Miller Parsons / Somerset, NJ SAMPLE ID: Sampling Method: Low flow purge (Monsoon Pump) **WELL PURGING** Static Water Level (TOC): 34.50 Depth to Well Bottom (TOC): 51.00 **CALCULATIONS:** Ft. of Water in Well X (GAL / FT) =Gallons 2-inch Casing: Ft. of Water in Wel Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 =Gallons 4-inch Casing: Ft. of Water in Well Gallons x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: Strong hydrocarbon odor Other: Clear, Sheen, NAPL blebs **FIELD TESTS PURGE PURGE PURGE** PURGE **PURGE PURGE PURGE PURGE** Time 1112 1127 1142 1117 1122 1132 1137 1147 Depth To Water (TOC) (ft) 36.51 37.00 39.00 38.20 38.00 39.05 38.50 37.90 Depth To Pump (TOC) (ft) 46.00 46.00 46.00 46.00 46.00 46.00 46.00 46.00 500 500 500 500 500 Flow Rate (ml/min) 500 500 500 Volume of Water Purged 0.75 2.25 5.25 1.5 3.0 3.75 4.5 6.0 6.94 7.02 7.03 7.04 7.04 7.08 7.09 7.10 pH (s.u.) Conductivity (mS/cm) 12.3 23.7 18.3 16.6 15.4 15.1 11.9 11.6 Turbidity (NTUs) 96.9 61.4 52.0 55.1 58.5 78.0 125 178 Dissolved Oxygen (mg/L) 1.34 0.92 0.78 0.69 0.68 0.67 0.66 0.65 Temperature (Degrees C) 16.05 16.14 16.11 16.34 16.37 16.81 16.90 17.10 -14 -17 -19 -21 -24 -27 -29 -32 ORP (mV) 7.1 13.2 10.7 9.6 8.9 8.6 8.0 6.5 Salinity (%) 13.6 11.2 10.2 9.5 9.2 8.9 8.2 7.8 TDS (g/L) **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's, TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1232; purged approximately 12.25 gallons. Flow cell cleaned out at 1227.

GROUNDWATER SAMPLING RECORD

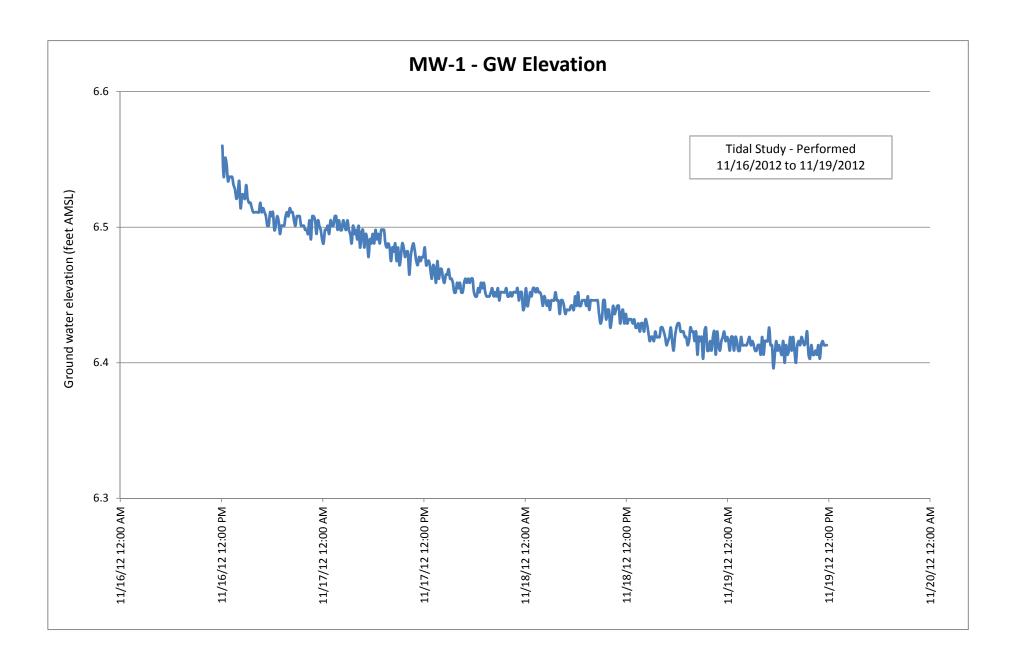
SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 4-8-13 Sampling Date: 4-8-13 Samplers: Parsons / Somerset, NJ Zohar Lavy SAMPLE ID: MW-7 Sampling Method: Low flow purge (Monsoon Pump) **WELL PURGING** Static Water Level (TOC): 34.50 Depth to Well Bottom (TOC): 51.00 **CALCULATIONS:** Ft. of Water in Well X (GAL / FT) =Gallons 2-inch Casing: Ft. of Water in Wel Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 =Gallons 4-inch Casing: Ft. of Water in Well Gallons x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION Odor: Strong hydrocarbon odor Other: Clear, Sheen, NAPL blebs **FIELD TESTS PURGE PURGE PURGE** PURGE **PURGE PURGE PURGE PURGE** Time 1152 1157 1202 1207 1217 1212 1222 1227 Depth To Water (TOC) (ft) 37.75 37.60 37.60 37.60 37.60 37.60 37.60 37.60 Depth To Pump (TOC) (ft) 46.00 46.00 46.00 46.00 46.00 46.00 46.00 46.00 500 500 500 500 500 500 Flow Rate (ml/min) 500 500 Volume of Water Purged 6.75 11.25 7.5 8.25 9.0 9.75 10.5 12.0 7.10 7.11 7.13 7.08 7.09 7.09 7.10 pH (s.u.) 7.11 Conductivity (mS/cm) 10.9 11.4 11.0 12.2 11.5 11.1 11.0 11.1 Turbidity (NTUs) 173 182 188 193 210 240 267 NA Dissolved Oxygen (mg/L) 0.63 0.61 0.60 0.59 0.59 0.58 0.62 0.68 Temperature (Degrees C) 17.20 17.00 16.70 16.47 16.50 16.53 16.57 16.51 -33 -34 -34 -34 -36 -37 -38 -37 ORP (mV) 6.3 5.8 6.0 6.5 6.4 6.3 6.1 Salinity (%) 6.3 6.7 7.2 6.4 6.8 7.1 6.9 6.8 6.9 TDS (g/L) **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's, TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1232; purged approximately 12.25 gallons. Flow cell cleaned out at 1227.

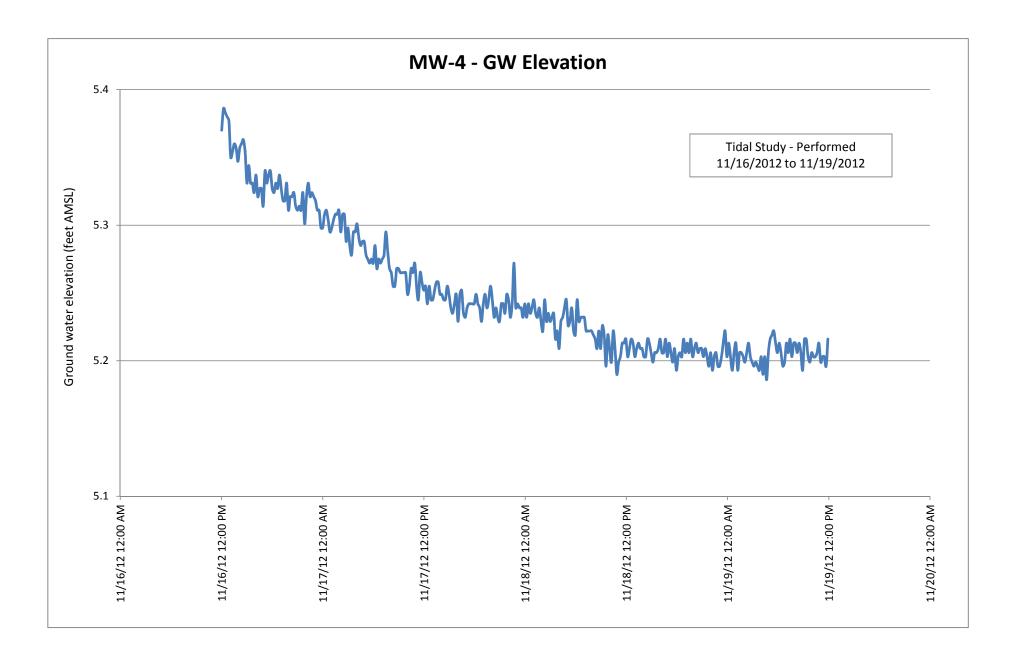
GROUNDWATER SAMPLING RECORD

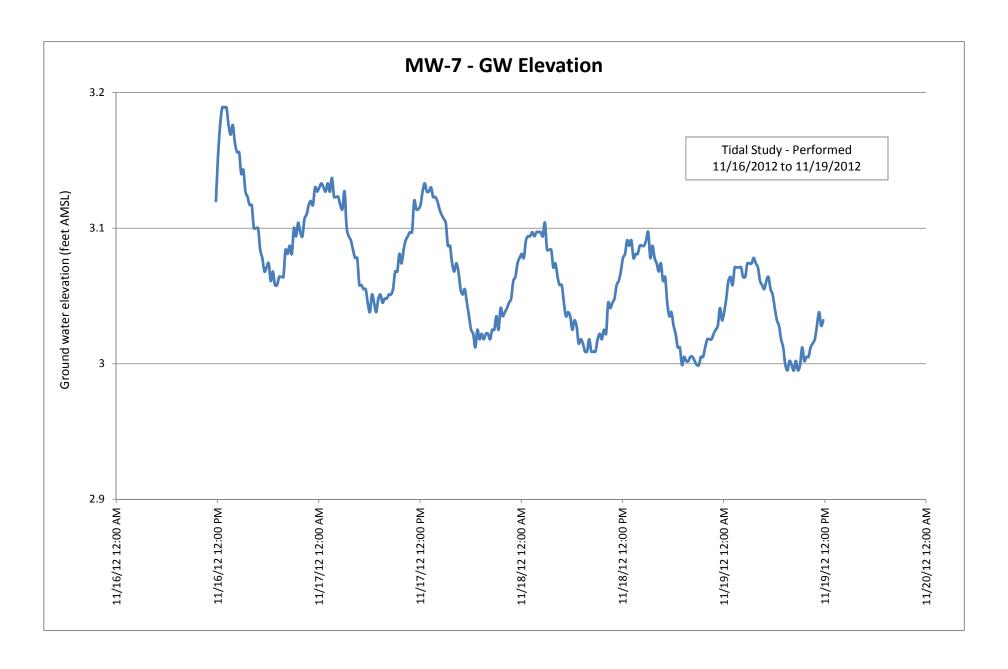
PROJECT NUMBER:	446110-04000			
Purge Date:	4-8-13			
Sampling Date:	4-8-13			
Samplers:	Zohar Lavy	of	Parsons / Somerset, NJ	
SAMPLE ID:	MW-7			
Sampling Method:	Low flow purge (Monsoon P	ump)		
ELL PURGING				
Static Water Level (TOC):	34.50			
Depth to Well Bottom (TOC):	51.00			
CALCULATIONS:	Ft. of Water in Well	X (GAL / FT) =	Gallons	
2-inch Casing:	Ft. of Water in Wel	x 0.16 =	Gallons	
3-inch Casing:	Ft. of Water in Well	x 0.32 =	Gallons	
4-inch Casing:	Ft. of Water in Well	x 0.64 =	Gallons	
Method:	Low Flow Pump			
SAMPLE DESCRIPTION				
Odor:	Strong hydrocarbon odor			
Other:	Clear, Sheen, NAPL blebs			
FIELD TESTS	Clear, Green, IVAL E Blebs			
ILLD ILGIO	SAMPLE			
Time	1232			
Depth To Water (TOC) (ft)	37.60			
Depth To Pump (TOC) (ft)	46.00			
Flow Rate (ml/min)	500			
Volume of Water Purged	12.25			
· ·	7.15			
pH (s.u.)	11.3			
Conductivity (mS/cm)				
Turbidity (NTUs)	17.9			
Dissolved Oxygen (mg/L)	0.73			
Temperature (Degrees C)	16.47			
ORP (mV)	-38			
Salinity (%)	6.4			
TDS (g/L)	7.0			
SAMPLE ANALYSIS / LABORAT	ORY			<u></u>
Analyze For:	TCL VOC's, TCL SVOCs, T	AL Metals, Cyanide		
Shipped Via:	Chemtech			
• • • • • • • • • • • • • • • • • • • •	Chemtech			
Laboratory				
Other Notes:	Sample collected at 1232; p	urged approximately 12.25 ga	illons. Flow cell cleaned out at 1227.	

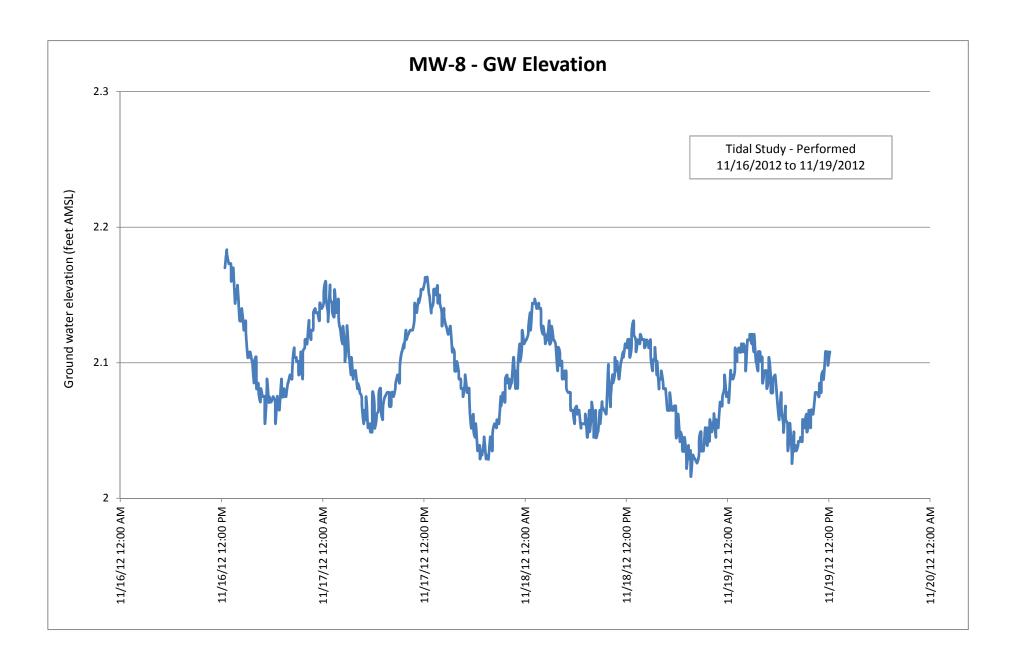
GROUNDWATER SAMPLING RECORD

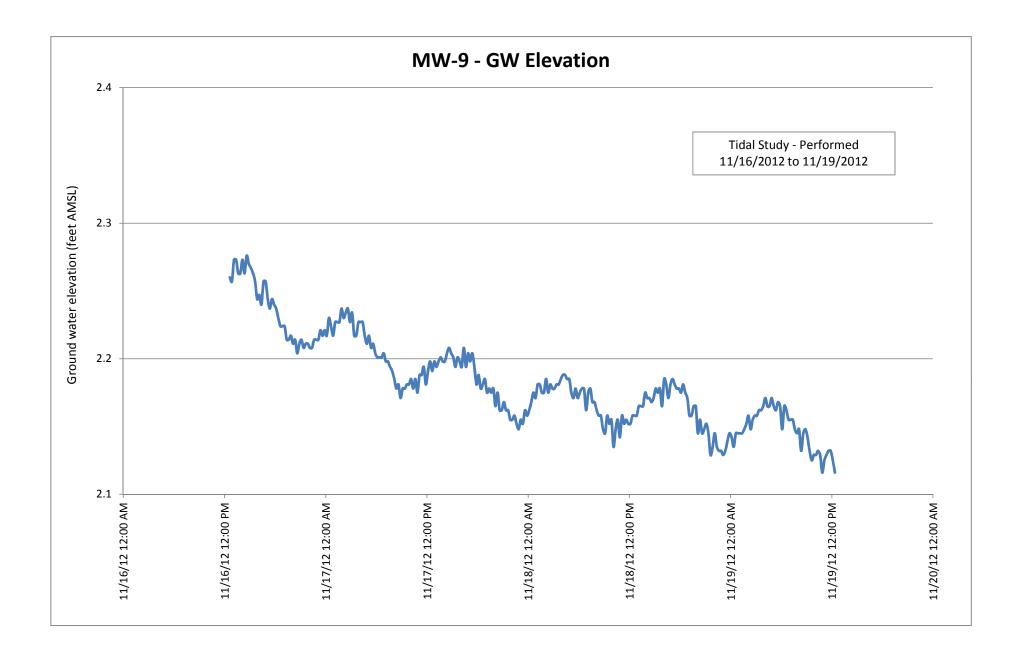
SITE NAME: Con Edison (Ludlow) PROJECT NUMBER: 446110-04000 Purge Date: 4-5-13 Sampling Date: 4-5-13 Samplers: Parsons / Somerset, NJ Zohar Lavy SAMPLE ID: MW-9 Sampling Method: Low flow purge (Monsoon Pump) **WELL PURGING** Static Water Level (TOC): 4.15 Depth to Well Bottom (TOC): 10.35 **CALCULATIONS:** Ft. of Water in Well X (GAL/FT) =Gallons 2-inch Casing: Ft. of Water in Wel Gallons x 0.16 =3-inch Casing: Ft. of Water in Well x 0.32 =Gallons 4-inch Casing: Ft. of Water in Well Gallons x 0.64 =Method: Low Flow Pump SAMPLE DESCRIPTION No Odor Odor: Other: Clear **FIELD TESTS PURGE PURGE PURGE** PURGE **PURGE PURGE** SAMPLE Time 1005 1010 1015 1020 1025 1030 1035 Depth To Water (TOC) (ft) 4.1 4.11 4.12 4.11 4.11 4.12 4.1 Depth To Pump (TOC) (ft) 9.35 9.35 9.35 9.35 9.35 9.35 9.35 ~300 Flow Rate (ml/min) ~350 ~350 ~350 ~300 ~300 ~300 Volume of Water Purged ~1.75 ~2.5 ~3.0 ~3.5 ~4.0 ~4.5 ~1 7.1 7.12 7.15 7.14 7.14 7.14 7.15 pH (s.u.) Conductivity (mS/cm) 6.56 6.88 6.55 6.35 6.53 6.61 6.35 Turbidity (NTUs) 82.9 22.5 2.0 0 0 0 0 Dissolved Oxygen (mg/L) 6.58 2.95 2.26 1.93 1.89 1.68 1.63 10.31 Temperature (Degrees C) 10.59 10.37 10.11 10.05 10.08 10.32 -51 -73 -77 -77 -82 -85 ORP (mV) -63 Salinity (%) 3.5 3.7 3.5 3.4 3.4 3.5 3.6 4.0 TDS (g/L) 4.14 4.34 4.12 4.0 4.14 4.15 **SAMPLE ANALYSIS / LABORATORY** Analyze For: TCL VOC's, TCL SVOCs, TAL Metals, Cyanide Shipped Via: Chemtech Laboratory Other Notes: Sample collected at 1035; purged approximately 4.5 gallons; began purging at 0950













11/8/2012 Mr. Zohar Lavy PARSONS GOVERNMENT SERVICES INC. 200 Cottontail Lane South

Somerset NJ 08873

Project Name: Ludlow

Project #:

Workorder #: 1210554

Dear Mr. Zohar Lavy

The following report includes the data for the above referenced project for sample(s) received on 10/25/2012 at Air Toxics Ltd.

The data and associated QC analyzed by Modified TO-15 are compliant with the project requirements or laboratory criteria with the exception of the deviations noted in the attached case narrative.

Thank you for choosing Air Toxics Ltd. for your air analysis needs. Air Toxics Ltd. is committed to providing accurate data of the highest quality. Please feel free to contact the Project Manager: Ausha Scott at 916-985-1000 if you have any questions regarding the data in this report.

Regards,

Ausha Scott

Project Manager



WORK ORDER #: 1210554

Work Order Summary

CLIENT: Mr. Zohar Lavy BILL TO: Accounts Payable-Syracuse

PARSONS GOVERNMENT PARSONS GOVERNMENT SERVICES

SERVICES INC. INC.

200 Cottontail Lane South 301 Plainfield Rd, Suite 350 Somerset, NJ 08873 Syracuse, NY 13212

PHONE: 732-537-3585 **P.O.** # 446110-02000

FAX: PROJECT # Ludlow

DATE RECEIVED: 10/25/2012 CONTACT: Ausha Scott 11/08/2012

			RECEIPT	FINAL
FRACTION #	NAME	<u>TEST</u>	VAC./PRES.	PRESSURE
01A	SV-3	Modified TO-15	6.5"Hg	5.0 psi
02A	SV-4	Modified TO-15	7.0"Hg	5.0 psi
03A	SV-1	Modified TO-15	7.5"Hg	5.0 psi
04A	Lab Blank	Modified TO-15	NA	NA
05A	CCV	Modified TO-15	NA	NA
06A	LCS	Modified TO-15	NA	NA
06AA	LCSD	Modified TO-15	NA	NA

	fleide flages	
CERTIFIED BY:		DATE: 11/08/12

Technical Director

Certfication numbers: AZ Licensure AZ0775, CA NELAP - 12282CA, NY NELAP - 11291, TX NELAP - T104704434-12-5, UT NELAP CA009332012-3, WA NELAP - C935

Name of Accrediting Agency: NELAP/ORELAP (Oregon Environmental Laboratory Accreditation Program) Accreditation number: CA300005, Effective date: 10/18/2011, Expiration date: 10/17/2012.

Eurofins Air Toxics Ltd. certifies that the test results contained in this report meet all requirements of the NELAC standards

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LABORATORY NARRATIVE EPA Method TO-15 PARSONS GOVERNMENT SERVICES INC. Workorder# 1210554

Three 6 Liter Summa Canister samples were received on October 25, 2012. The laboratory performed analysis via EPA Method TO-15 using GC/MS in the full scan mode.

This workorder was independently validated prior to submittal using 'USEPA National Functional Guidelines' as generally applied to the analysis of volatile organic compounds in air. A rules-based, logic driven, independent validation engine was employed to assess completeness, evaluate pass/fail of relevant project quality control requirements and verification of all quantified amounts.

Receiving Notes

There were no receiving discrepancies.

Analytical Notes

All Quality Control Limit exceedances and affected sample results are noted by flags. Each flag is defined at the bottom of this Case Narrative and on each Sample Result Summary page.

Definition of Data Qualifying Flags

Eight qualifiers may have been used on the data analysis sheets and indicates as follows:

- B Compound present in laboratory blank greater than reporting limit (background subtraction not performed).
 - J Estimated value.
 - E Exceeds instrument calibration range.
 - S Saturated peak.
 - Q Exceeds quality control limits.
 - U Compound analyzed for but not detected above the reporting limit.
 - UJ- Non-detected compound associated with low bias in the CCV and/or LCS.
 - N The identification is based on presumptive evidence.

File extensions may have been used on the data analysis sheets and indicates as follows:

a-File was requantified

b-File was quantified by a second column and detector

r1-File was requantified for the purpose of reissue



Summary of Detected Compounds EPA METHOD TO-15 GC/MS FULL SCAN

Client Sample ID: SV-3 Lab ID#: 1210554-01A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)	
Acetone	8.6	64	20	150	
Toluene	0.86	6.2	3.2	23	
Tetrachloroethene	0.86	15	5.8	100	
m,p-Xylene	0.86	2.6	3.7	11	

Client Sample ID: SV-4 Lab ID#: 1210554-02A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,3-Butadiene	0.88	2.0	1.9	4.4
Acetone	8.8	91	21	220
Carbon Disulfide	3.5	4.5	11	14
Hexane	0.88	100	3.1	360
Benzene	0.88	2.7	2.8	8.6
Heptane	0.88	47	3.6	190
Toluene	0.88	16	3.3	60
Tetrachloroethene	0.88	3.6	5.9	24
Ethyl Benzene	0.88	1.3	3.8	5.8
m,p-Xylene	0.88	4.6	3.8	20
o-Xylene	0.88	1.6	3.8	7.0

Client Sample ID: SV-1

Lab ID#: 1210554-03A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Freon 12	0.90	1.2	4.4	5.9
1,3-Butadiene	0.90	15	2.0	34
Acetone	9.0	110	21	260
Carbon Disulfide	3.6	5.4	11	17
Hexane	0.90	7.6	3.2	27
2-Butanone (Methyl Ethyl Ketone)	3.6	3.8	10	11
1,1,1-Trichloroethane	0.90	1.2	4.9	6.5
Cyclohexane	0.90	2.3	3.1	7.9



Summary of Detected Compounds EPA METHOD TO-15 GC/MS FULL SCAN

Client Sample 1	ID: SV-1
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Lab ID#: 1210554-03A				
2,2,4-Trimethylpentane	0.90	2.5	4.2	12
Benzene	0.90	6.2	2.8	20
Heptane	0.90	3.2	3.7	13
Toluene	0.90	9.2	3.4	35
Tetrachloroethene	0.90	4.2	6.1	28
Ethyl Benzene	0.90	1.2	3.9	5.4
m,p-Xylene	0.90	4.4	3.9	19
o-Xylene	0.90	1.2	3.9	5.2



Client Sample ID: SV-3 Lab ID#: 1210554-01A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name:	p102639	Date of Collection: 10/24/12 11:55:00 A
Dil. Factor:	1.71	Date of Analysis: 10/27/12 04:55 PM

Dil. Factor:	1.71	Date of Analysis: 10/27/12 04:55 PM		
	Rpt. Limit	Amount	Rpt. Limit	Amount
Compound	(ppbv)	(ppbv)	(ug/m3)	(ug/m3)
Freon 12	0.86	Not Detected	4.2	Not Detected
Freon 114	0.86	Not Detected	6.0	Not Detected
Chloromethane	8.6	Not Detected	18	Not Detected
Vinyl Chloride	0.86	Not Detected	2.2	Not Detected
1,3-Butadiene	0.86	Not Detected	1.9	Not Detected
Bromomethane	8.6	Not Detected	33	Not Detected
Chloroethane	3.4	Not Detected	9.0	Not Detected
Freon 11	0.86	Not Detected	4.8	Not Detected
Ethanol	3.4	Not Detected	6.4	Not Detected
Freon 113	0.86	Not Detected	6.6	Not Detected
1,1-Dichloroethene	0.86	Not Detected	3.4	Not Detected
Acetone	8.6	64	20	150
2-Propanol	3.4	Not Detected	8.4	Not Detected
Carbon Disulfide	3.4	Not Detected	11	Not Detected
3-Chloropropene	3.4	Not Detected	11	Not Detected
Methylene Chloride	8.6	Not Detected	30	Not Detected
Methyl tert-butyl ether	0.86	Not Detected	3.1	Not Detected
trans-1,2-Dichloroethene	0.86	Not Detected	3.4	Not Detected
Hexane	0.86	Not Detected	3.0	Not Detected
1,1-Dichloroethane	0.86	Not Detected	3.5	Not Detected
2-Butanone (Methyl Ethyl Ketone)	3.4	Not Detected	10	Not Detected
cis-1,2-Dichloroethene	0.86	Not Detected	3.4	Not Detected
Tetrahydrofuran	0.86	Not Detected	2.5	Not Detected
Chloroform	0.86	Not Detected	4.2	Not Detected
1,1,1-Trichloroethane	0.86	Not Detected	4.7	Not Detected
Cyclohexane	0.86	Not Detected	2.9	Not Detected
Carbon Tetrachloride	0.86	Not Detected	5.4	Not Detected
2,2,4-Trimethylpentane	0.86	Not Detected	4.0	Not Detected
Benzene	0.86	Not Detected	2.7	Not Detected
1,2-Dichloroethane	0.86	Not Detected	3.5	Not Detected
Heptane	0.86	Not Detected	3.5	Not Detected
Trichloroethene	0.86	Not Detected	4.6	Not Detected
1,2-Dichloropropane	0.86	Not Detected	4.0	Not Detected
1,4-Dioxane	3.4	Not Detected	12	Not Detected
Bromodichloromethane	0.86	Not Detected	5.7	Not Detected
cis-1,3-Dichloropropene	0.86	Not Detected	3.9	Not Detected
4-Methyl-2-pentanone	0.86	Not Detected	3.5	Not Detected
Toluene	0.86	6.2	3.2	23
trans-1,3-Dichloropropene	0.86	Not Detected	3.9	Not Detected
1,1,2-Trichloroethane	0.86	Not Detected	4.7	Not Detected
Tetrachloroethene	0.86	15	5.8	100
2-Hexanone	3.4	Not Detected	14	Not Detected



Client Sample ID: SV-3 Lab ID#: 1210554-01A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name:	p102639	Date of Collection: 10/24/12 11:55:00 A
Dil. Factor:	1.71	Date of Analysis: 10/27/12 04:55 PM

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Dibromochloromethane	0.86	Not Detected	7.3	Not Detected
1,2-Dibromoethane (EDB)	0.86	Not Detected	6.6	Not Detected
Chlorobenzene	0.86	Not Detected	3.9	Not Detected
Ethyl Benzene	0.86	Not Detected	3.7	Not Detected
m,p-Xylene	0.86	2.6	3.7	11
o-Xylene	0.86	Not Detected	3.7	Not Detected
Styrene	0.86	Not Detected	3.6	Not Detected
Bromoform	0.86	Not Detected	8.8	Not Detected
Cumene	0.86	Not Detected	4.2	Not Detected
1,1,2,2-Tetrachloroethane	0.86	Not Detected	5.9	Not Detected
Propylbenzene	0.86	Not Detected	4.2	Not Detected
4-Ethyltoluene	0.86	Not Detected	4.2	Not Detected
1,3,5-Trimethylbenzene	0.86	Not Detected	4.2	Not Detected
1,2,4-Trimethylbenzene	0.86	Not Detected	4.2	Not Detected
1,3-Dichlorobenzene	0.86	Not Detected	5.1	Not Detected
1,4-Dichlorobenzene	0.86	Not Detected	5.1	Not Detected
alpha-Chlorotoluene	0.86	Not Detected	4.4	Not Detected
1,2-Dichlorobenzene	0.86	Not Detected	5.1	Not Detected
1,2,4-Trichlorobenzene	3.4	Not Detected	25	Not Detected
Hexachlorobutadiene	3.4	Not Detected	36	Not Detected

Container Type: 6 Liter Summa Canister

Surrogates	%Recovery	Method Limits
Toluene-d8	96	70-130
1,2-Dichloroethane-d4	99	70-130
4-Bromofluorobenzene	103	70-130



Client Sample ID: SV-4 Lab ID#: 1210554-02A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: p102640 Date of Collection: 10/24/12 12:10:00 P
Dil. Factor: 1.75 Date of Analysis: 10/27/12 05:38 PM

Dil. Factor:	1.75	Date of Analysis: 10/27/12 05:38		
	Rpt. Limit	Amount	Rpt. Limit	Amount
Compound	(ppbv)	(ppbv)	(ug/m3)	(ug/m3)
Freon 12	0.88	Not Detected	4.3	Not Detected
Freon 114	0.88	Not Detected	6.1	Not Detected
Chloromethane	8.8	Not Detected	18	Not Detected
Vinyl Chloride	0.88	Not Detected	2.2	Not Detected
1,3-Butadiene	0.88	2.0	1.9	4.4
Bromomethane	8.8	Not Detected	34	Not Detected
Chloroethane	3.5	Not Detected	9.2	Not Detected
Freon 11	0.88	Not Detected	4.9	Not Detected
Ethanol	3.5	Not Detected	6.6	Not Detected
Freon 113	0.88	Not Detected	6.7	Not Detected
1,1-Dichloroethene	0.88	Not Detected	3.5	Not Detected
Acetone	8.8	91	21	220
2-Propanol	3.5	Not Detected	8.6	Not Detected
Carbon Disulfide	3.5	4.5	11	14
3-Chloropropene	3.5	Not Detected	11	Not Detected
Methylene Chloride	8.8	Not Detected	30	Not Detected
Methyl tert-butyl ether	0.88	Not Detected	3.2	Not Detected
trans-1,2-Dichloroethene	0.88	Not Detected	3.5	Not Detected
Hexane	0.88	100	3.1	360
1,1-Dichloroethane	0.88	Not Detected	3.5	Not Detected
2-Butanone (Methyl Ethyl Ketone)	3.5	Not Detected	10	Not Detected
cis-1,2-Dichloroethene	0.88	Not Detected	3.5	Not Detected
Tetrahydrofuran	0.88	Not Detected	2.6	Not Detected
Chloroform	0.88	Not Detected	4.3	Not Detected
1,1,1-Trichloroethane	0.88	Not Detected	4.8	Not Detected
Cyclohexane	0.88	Not Detected	3.0	Not Detected
Carbon Tetrachloride	0.88	Not Detected	5.5	Not Detected
2,2,4-Trimethylpentane	0.88	Not Detected	4.1	Not Detected
Benzene	0.88	2.7	2.8	8.6
1,2-Dichloroethane	0.88	Not Detected	3.5	Not Detected
Heptane	0.88	47	3.6	190
Trichloroethene	0.88	Not Detected	4.7	Not Detected
1,2-Dichloropropane	0.88	Not Detected	4.0	Not Detected
1,4-Dioxane	3.5	Not Detected	13	Not Detected
Bromodichloromethane	0.88	Not Detected	5.9	Not Detected
cis-1,3-Dichloropropene	0.88	Not Detected	4.0	Not Detected
4-Methyl-2-pentanone	0.88	Not Detected	3.6	Not Detected
Toluene	0.88	16	3.3	60
trans-1,3-Dichloropropene	0.88	Not Detected	4.0	Not Detected
1,1,2-Trichloroethane	0.88	Not Detected	4.8	Not Detected
Tetrachloroethene	0.88	3.6	5.9	24
2-Hexanone	3.5	Not Detected	14	Not Detected



Client Sample ID: SV-4 Lab ID#: 1210554-02A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name:	p102640	Date of Collection: 10/24/12 12:10:00 P
Dil. Factor:	1.75	Date of Analysis: 10/27/12 05:38 PM

2 at 0 0 7 m lang 0 10 1	0,=1,1=00.00
unt Rpt. Limit ov) (ug/m3)	Amount (ug/m3)
ected 7.4	Not Detected
ected 6.7	Not Detected
ected 4.0	Not Detected
3.8	5.8
3.8	20
3.8	7.0
ected 3.7	Not Detected
ected 9.0	Not Detected
ected 4.3	Not Detected
ected 6.0	Not Detected
ected 4.3	Not Detected
ected 5.3	Not Detected
ected 5.3	Not Detected
ected 4.5	Not Detected
ected 5.3	Not Detected
ected 26	Not Detected
	Not Detected
ttttt	tected 9.0 tected 4.3 tected 6.0 tected 4.3 tected 4.3 tected 4.3 tected 4.3 tected 5.3

Container Type: 6 Liter Summa Canister

Surrogates	%Recovery	Method Limits
Toluene-d8	101	70-130
1,2-Dichloroethane-d4	101	70-130
4-Bromofluorobenzene	97	70-130



Client Sample ID: SV-1 Lab ID#: 1210554-03A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name:	p102641	Date of Collection: 10/24/12 2:28:00 PM
Dil. Factor:	1.79	Date of Analysis: 10/27/12 06:09 PM

Compound Rpt. Limit (ppbv) Amount (ppbv) Rpt. Limit (ug/m3) Amount (ug/m3) Freon 12 0.90 1.2 4.4 5.9 Freon 114 0.90 Not Detected 6.2 Not Detected Chloromethane 9.0 Not Detected 18 Not Detected Linyl Chloride 0.90 Not Detected 2.3 Not Detected 1,3-Butadiene 9.0 Not Detected 3.5 Not Detected Chloromethane 3.6 Not Detected 9.4 Not Detected Chloroethane 3.6 Not Detected 5.0 Not Detected Chloroethane 3.6 Not Detected 6.7 Not Detected Freon 11 9.0 Not Detected 6.8 Not Detected Freon 113 0.90 Not Detected 6.8 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected 4,1-Dichloroethene 0.90 Not Detected 3.8 Not Detected 2-Propanol 3.6	Dil. Factor:	1.79	Date	of Analysis: 10/2	7/12 06:09 PM
Freon 12 0.90 1.2 4.4 5.9 Freon 114 0.90 Not Detected 6.2 Not Detected Chloromethane 9.0 Not Detected 18 Not Detected Vinyl Chloride 0.90 Not Detected 2.3 Not Detected 1,3-Butadiene 0.90 15 2.0 34 Brommethane 9.0 Not Detected 3.5 Not Detected Chloroethane 3.6 Not Detected 3.0 Not Detected Freon 11 0.90 Not Detected 5.0 Not Detected Freon 13 0.90 Not Detected 6.7 Not Detected Freon 13 0.90 Not Detected 6.8 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected 1,1-Dichloroethene 0.90 Not Detected 8.8 Not Detected 2-Propanol 3.6 Not Detected 8.8 Not Detected 4-Propanol 3.6 Not Detected 1.1		Rpt. Limit	Amount	Rpt. Limit	Amount
Freon 114 0.90 Not Detected 6.2 Not Detected Chloromethane 9.0 Not Detected 18 Not Detected Unity Chloride 0.90 Not Detected 2.3 Not Detected 1,3-Butadiene 0.90 15 2.0 34 Bromomethane 9.0 Not Detected 3.5 Not Detected Chloroethane 3.6 Not Detected 5.0 Not Detected Freon 11 0.90 Not Detected 6.7 Not Detected Ethanol 3.6 Not Detected 6.8 Not Detected Freon 113 0.90 Not Detected 3.5 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected 4,1-Dichloroethene 9.0 Not Detected 3.8 Not Detected 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 Not Detected 1.1 17 3-Chiloropropropen 3.6 Not Detected	Compound	(ppbv)	(ppbv)	(ug/m3)	(ug/m3)
Chloromethane 9.0 Not Detected 18 Not Detected Vinyl Chloride 0.90 Not Detected 2.3 Not Detected 1.3-Butadiene 0.90 15 2.0 34 Bromomethane 9.0 Not Detected 35 Not Detected Chloroethane 3.6 Not Detected 5.0 Not Detected Freon 11 0.90 Not Detected 6.7 Not Detected Ethanol 3.6 Not Detected 6.7 Not Detected Freon 113 0.90 Not Detected 3.5 Not Detected 1,-Dichloroethene 0.90 Not Detected 3.5 Not Detected Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 3.8 Not Detected Acetone 3.0 Not Detected 3.8 Not Detected Methyler Chloride 3.6 Not Detected 3.1 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2	Freon 12	0.90	1.2	4.4	5.9
Vinyl Chloride 0.90 Not Detected 2.3 Not Detected 1,3-Butadiene 0.90 15 2.0 34 Brommethane 9.0 Not Detected 35 Not Detected Chloroethane 3.6 Not Detected 9.4 Not Detected Freon 11 0.90 Not Detected 6.7 Not Detected Ethanol 3.6 Not Detected 6.8 Not Detected Freon 113 0.90 Not Detected 3.5 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 Not Detected 4.8 Not Detected Achtoryporpopene 3.6 Not Detected 3.1 Not Detected Methyl ten-butyl ether 0.90 Not Detected 3.1 Not Detected Methyl ten-butyl ether 0.90 Not Detected	Freon 114	0.90	Not Detected	6.2	Not Detected
1,3-Butadiene 0.90 15 2.0 34 Bromomethane 9.0 Not Detected 35 Not Detected Chloroethane 3.6 Not Detected 9.0 Not Detected Freon 11 0.90 Not Detected 5.0 Not Detected Ethanol 3.6 Not Detected 6.8 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 Not Detected 11 Not Detected Carbon Chloride 9.0 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 3.1 Not Detected Methylene Chloride 9.0 Not Detected 3.2 Not Detected Methylene Chloride 9.0 Not Detected 3.2 Not Detected Methylene Chloride 9.0 Not Detected	Chloromethane	9.0	Not Detected	18	Not Detected
Bromomethane	Vinyl Chloride	0.90	Not Detected	2.3	Not Detected
Chloroethane 3.6 Not Detected 9.4 Not Detected Freon 11 0.90 Not Detected 5.0 Not Detected Ethanol 3.6 Not Detected 6.7 Not Detected Freon 113 0.90 Not Detected 6.8 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 3.1 Not Detected Methylene Chloride 9.0 Not Detected 3.2 Not Detected Methylene Chloride 9.0 Not Detected 3.5 Not Detected Methylene Chloride 9.0 Not Detected 3.5 Not Detected Hexane 0.90 Not Detecte	1,3-Butadiene	0.90	15	2.0	34
Freon 11 0.90 Not Detected 5.0 Not Detected Ethanol 3.6 Not Detected 6.7 Not Detected Freon 113 0.90 Not Detected 6.8 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 5.4 11 17 3-Chloropropene 3.6 Not Detected 11 Not Detected Methyler-buryle ether 0.90 Not Detected 3.1 Not Detected Methyler-buryle ether 0.90 Not Detected 3.2 Not Detected trans-1,2-Dichloroethene 0.90 Not Detected 3.2 Not Detected trans-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected trans-1,2-Dichloroethane 0.90 Not Detected 3.5 Not Detected trans-1,2-Dichloroethane 0.90 Not Detected 3.5 Not Detected trans-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected trans-1,2-Dichloroethene 0.90 Not Detected 3.5 <	Bromomethane	9.0	Not Detected	35	Not Detected
Ethanol 3.6 Not Detected 6.7 Not Detected Freon 113 0.90 Not Detected 6.8 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 31 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.5 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.5 Not Detected 1,2-Di	Chloroethane	3.6	Not Detected	9.4	Not Detected
Freon 113 0.90 Not Detected 6.8 Not Detected 1,1-Dichloroethene 0.90 Not Detected 3.5 Not Detected Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 5.4 11 17 3-Chloropropene 3.6 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 31 Not Detected Methylene Chloride 9.0 Not Detected 3.2 Not Detected Methylene Chloride 9.0 Not Detected 3.2 Not Detected Methylene Chloride 9.0 Not Detected 3.2 Not Detected trans-1,2-Dichloroethene 0.90 Not Detected 3.6 Not Detected 4-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Tetrahydrofuran 0.90	Freon 11	0.90	Not Detected	5.0	Not Detected
1,1-Dichloroethene	Ethanol	3.6	Not Detected	6.7	Not Detected
Acetone 9.0 110 21 260 2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 5.4 11 17 3-Chloropropene 3.6 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 31 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Hexane 0.90 Not Detected 3.5 Not Detected Hexane 0.90 7.6 3.2 2.7 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethane 0.90 Not Detected 3.5 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,1-Tichloroethane 0.90 Not Detected	Freon 113	0.90	Not Detected	6.8	Not Detected
2-Propanol 3.6 Not Detected 8.8 Not Detected Carbon Disulfide 3.6 5.4 11 17 3-Chloropropene 3.6 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 31 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Hexane 0.90 7.6 3.2 27 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 3.5 Not Detected Chioroform 0.90 Not Detected 4.4 Not Detected Chioroform 0.90 Not Detected 4.4 Not Detected Chioroform 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1	1,1-Dichloroethene	0.90	Not Detected	3.5	Not Detected
Carbon Disulfide 3.6 5.4 11 17 3-Chloropropene 3.6 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 31 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected trans-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Hexane 0.90 7.6 3.2 27 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected Chloroform 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Det		9.0	110	21	260
3-Chloropropene 3.6 Not Detected 11 Not Detected Methylene Chloride 9.0 Not Detected 31 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected trans-1,2-Dichloroethene 0.90 7.6 3.2 27 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 3.5 Not Detected Chloroform 0.90 Not Detected 3.6 Not Detected 1,1,1-Trichloroethane 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 Not Detected 4.4 Not Detected 2,2,4-Trimethylpentane 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 0.90 3.2 3.7 13	2-Propanol	3.6	Not Detected	8.8	Not Detected
Methylene Chloride 9.0 Not Detected 31 Not Detected Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected Itanal Project 0.90 Not Detected 3.5 Not Detected Hexane 0.90 7.6 3.2 27 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 3.5 Not Detected Chloroform 0.90 Not Detected 2.6 Not Detected 1,1-Trichloroethane 0.90 Not Detected 4.4 Not Detected Cyclohexane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 Not Detected 5.6 Not Detected 2,2-Dichloroptethane 0.90	Carbon Disulfide	3.6	5.4	11	17
Methyl tert-butyl ether 0.90 Not Detected 3.2 Not Detected trans-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Hexane 0.90 7.6 3.2 27 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected Chloroform 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Cyclohexane 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 Not Detected 3.6 <td>3-Chloropropene</td> <td>3.6</td> <td>Not Detected</td> <td>11</td> <td>Not Detected</td>	3-Chloropropene	3.6	Not Detected	11	Not Detected
trans-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Hexane 0.90 7.6 3.2 27 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8	Methylene Chloride	9.0	Not Detected	31	Not Detected
Hexane 0.90 7.6 3.2 27 1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 8.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected	Methyl tert-butyl ether	0.90	Not Detected	3.2	Not Detected
1,1-Dichloroethane 0.90 Not Detected 3.6 Not Detected 2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,2-Dichloropropane 0.90 Not Detected	trans-1,2-Dichloroethene	0.90	Not Detected	3.5	Not Detected
2-Butanone (Methyl Ethyl Ketone) 3.6 3.8 10 11 cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected	Hexane	0.90	7.6	3.2	27
cis-1,2-Dichloroethene 0.90 Not Detected 3.5 Not Detected Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 4.1 Not Detected Bromodichloromethane 0.90 Not Detected <t< td=""><td>1,1-Dichloroethane</td><td>0.90</td><td>Not Detected</td><td>3.6</td><td>Not Detected</td></t<>	1,1-Dichloroethane	0.90	Not Detected	3.6	Not Detected
Tetrahydrofuran 0.90 Not Detected 2.6 Not Detected Chloroform 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 3.2 3.7 13 Trichloroethane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 4.1 Not Detected Bromodichloromethane 0.90 Not Detected 4.1 Not Detected cis-1,3-Dichloropropene 0.90 Not Detected 4.1	2-Butanone (Methyl Ethyl Ketone)	3.6	3.8	10	11
Chloroform 0.90 Not Detected 4.4 Not Detected 1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloroethene 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 4.1 Not Detected Bromodichloromethane 0.90 Not Detected 4.1 Not Detected cis-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 4-Methyl-2-pentanone 0.90 Not Detecte	cis-1,2-Dichloroethene	0.90	Not Detected	3.5	Not Detected
1,1,1-Trichloroethane 0.90 1.2 4.9 6.5 Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 13 Not Detected 1,4-Dioxane 3.6 Not Detected 13 Not Detected Bromodichloromethane 0.90 Not Detected 4.1 Not Detected 4-Methyl-2-pentanone 0.90 Not Detected 3.7 Not Detected Toluene 0.90 Not Detected 4.1 <td>Tetrahydrofuran</td> <td>0.90</td> <td>Not Detected</td> <td>2.6</td> <td>Not Detected</td>	Tetrahydrofuran	0.90	Not Detected	2.6	Not Detected
Cyclohexane 0.90 2.3 3.1 7.9 Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 13 Not Detected Bromodichloromethane 0.90 Not Detected 6.0 Not Detected cis-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 4-Methyl-2-pentanone 0.90 Not Detected 3.7 Not Detected Toluene 0.90 Not Detected 4.1 Not Detected 1,1,2-Trichloroethane 0.90 Not D	Chloroform	0.90	Not Detected	4.4	Not Detected
Carbon Tetrachloride 0.90 Not Detected 5.6 Not Detected 2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 13 Not Detected Bromodichloromethane 0.90 Not Detected 6.0 Not Detected cis-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 4-Methyl-2-pentanone 0.90 Not Detected 3.7 Not Detected Toluene 0.90 9.2 3.4 35 trans-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 1,1,2-Trichloroethane 0.90 Not Detected 4.9 Not Detected Tetrachloroethene 0.90 <td>1,1,1-Trichloroethane</td> <td>0.90</td> <td>1.2</td> <td>4.9</td> <td>6.5</td>	1,1,1-Trichloroethane	0.90	1.2	4.9	6.5
2,2,4-Trimethylpentane 0.90 2.5 4.2 12 Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 3.2 3.7 13 Trichloroethene 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 13 Not Detected Bromodichloromethane 0.90 Not Detected 6.0 Not Detected cis-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 4-Methyl-2-pentanone 0.90 Not Detected 3.7 Not Detected Tolluene 0.90 9.2 3.4 35 trans-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 1,1,2-Trichloroethane 0.90 Not Detected 4.9 Not Detected Tetrachloroethene 0.90 4.2	Cyclohexane	0.90	2.3	3.1	7.9
Benzene 0.90 6.2 2.8 20 1,2-Dichloroethane 0.90 Not Detected 3.6 Not Detected Heptane 0.90 3.2 3.7 13 Trichloroethene 0.90 Not Detected 4.8 Not Detected 1,2-Dichloropropane 0.90 Not Detected 4.1 Not Detected 1,4-Dioxane 3.6 Not Detected 13 Not Detected Bromodichloromethane 0.90 Not Detected 6.0 Not Detected cis-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 4-Methyl-2-pentanone 0.90 Not Detected 3.7 Not Detected Toluene 0.90 9.2 3.4 35 trans-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 1,1,2-Trichloroethane 0.90 Not Detected 4.9 Not Detected Tetrachloroethene 0.90 4.2 6.1 28	Carbon Tetrachloride	0.90	Not Detected	5.6	Not Detected
1,2-Dichloroethane0.90Not Detected3.6Not DetectedHeptane0.903.23.713Trichloroethene0.90Not Detected4.8Not Detected1,2-Dichloropropane0.90Not Detected4.1Not Detected1,4-Dioxane3.6Not Detected13Not DetectedBromodichloromethane0.90Not Detected6.0Not Detectedcis-1,3-Dichloropropene0.90Not Detected4.1Not Detected4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	2,2,4-Trimethylpentane	0.90		4.2	
Heptane0.903.23.713Trichloroethene0.90Not Detected4.8Not Detected1,2-Dichloropropane0.90Not Detected4.1Not Detected1,4-Dioxane3.6Not Detected13Not DetectedBromodichloromethane0.90Not Detected6.0Not Detectedcis-1,3-Dichloropropene0.90Not Detected4.1Not Detected4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	Benzene	0.90	6.2	2.8	20
Trichloroethene0.90Not Detected4.8Not Detected1,2-Dichloropropane0.90Not Detected4.1Not Detected1,4-Dioxane3.6Not Detected13Not DetectedBromodichloromethane0.90Not Detected6.0Not Detectedcis-1,3-Dichloropropene0.90Not Detected4.1Not Detected4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	1,2-Dichloroethane	0.90	Not Detected	3.6	Not Detected
1,2-Dichloropropane0.90Not Detected4.1Not Detected1,4-Dioxane3.6Not Detected13Not DetectedBromodichloromethane0.90Not Detected6.0Not Detectedcis-1,3-Dichloropropene0.90Not Detected4.1Not Detected4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	Heptane	0.90	3.2	3.7	13
1,4-Dioxane3.6Not Detected13Not DetectedBromodichloromethane0.90Not Detected6.0Not Detectedcis-1,3-Dichloropropene0.90Not Detected4.1Not Detected4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	Trichloroethene	0.90	Not Detected	4.8	Not Detected
Bromodichloromethane0.90Not Detected6.0Not Detectedcis-1,3-Dichloropropene0.90Not Detected4.1Not Detected4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	1,2-Dichloropropane				
cis-1,3-Dichloropropene0.90Not Detected4.1Not Detected4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	1,4-Dioxane		Not Detected	13	Not Detected
4-Methyl-2-pentanone0.90Not Detected3.7Not DetectedToluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	Bromodichloromethane	0.90	Not Detected	6.0	Not Detected
Toluene0.909.23.435trans-1,3-Dichloropropene0.90Not Detected4.1Not Detected1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	cis-1,3-Dichloropropene	0.90	Not Detected	4.1	Not Detected
trans-1,3-Dichloropropene 0.90 Not Detected 4.1 Not Detected 1,1,2-Trichloroethane 0.90 Not Detected 4.9 Not Detected Tetrachloroethene 0.90 4.2 6.1 28	4-Methyl-2-pentanone	0.90	Not Detected	3.7	Not Detected
1,1,2-Trichloroethane0.90Not Detected4.9Not DetectedTetrachloroethene0.904.26.128	Toluene	0.90	9.2	3.4	35
Tetrachloroethene 0.90 4.2 6.1 28	trans-1,3-Dichloropropene	0.90	Not Detected	4.1	Not Detected
	1,1,2-Trichloroethane	0.90	Not Detected	4.9	Not Detected
2-Hexanone 3.6 Not Detected 15 Not Detected	Tetrachloroethene	0.90	4.2	6.1	28
	2-Hexanone	3.6	Not Detected	15	Not Detected



Client Sample ID: SV-1 Lab ID#: 1210554-03A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name:	p102641	Date of Collection: 10/24/12 2:28:00 PM
Dil. Factor:	1.79	Date of Analysis: 10/27/12 06:09 PM

J 1 4.01011	1170	Dute	or Amaryolo. 10/2	7712 00:00 T III
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Dibromochloromethane	0.90	Not Detected	7.6	Not Detected
1,2-Dibromoethane (EDB)	0.90	Not Detected	6.9	Not Detected
Chlorobenzene	0.90	Not Detected	4.1	Not Detected
Ethyl Benzene	0.90	1.2	3.9	5.4
m,p-Xylene	0.90	4.4	3.9	19
o-Xylene	0.90	1.2	3.9	5.2
Styrene	0.90	Not Detected	3.8	Not Detected
Bromoform	0.90	Not Detected	9.2	Not Detected
Cumene	0.90	Not Detected	4.4	Not Detected
1,1,2,2-Tetrachloroethane	0.90	Not Detected	6.1	Not Detected
Propylbenzene	0.90	Not Detected	4.4	Not Detected
4-Ethyltoluene	0.90	Not Detected	4.4	Not Detected
1,3,5-Trimethylbenzene	0.90	Not Detected	4.4	Not Detected
1,2,4-Trimethylbenzene	0.90	Not Detected	4.4	Not Detected
1,3-Dichlorobenzene	0.90	Not Detected	5.4	Not Detected
1,4-Dichlorobenzene	0.90	Not Detected	5.4	Not Detected
alpha-Chlorotoluene	0.90	Not Detected	4.6	Not Detected
1,2-Dichlorobenzene	0.90	Not Detected	5.4	Not Detected
1,2,4-Trichlorobenzene	3.6	Not Detected	26	Not Detected
Hexachlorobutadiene	3.6	Not Detected	38	Not Detected

Container Type: 6 Liter Summa Canister

		Method Limits	
Surrogates	%Recovery		
Toluene-d8	96	70-130	
1,2-Dichloroethane-d4	97	70-130	
4-Bromofluorobenzene	101	70-130	



Client Sample ID: Lab Blank Lab ID#: 1210554-04A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name:	p102623	Date of Collection: NA
Dil. Factor:	1.00	Date of Analysis: 10/26/12 10:44 PM

Dil. Factor: 1.00 Date of Analysis: 1		of Analysis: 10/2	0/26/12 10:44 PM	
	Rpt. Limit	Amount	Rpt. Limit	Amount
Compound	(ppbv)	(ppbv)	(ug/m3)	(ug/m3)
Freon 12	0.50	Not Detected	2.5	Not Detected
Freon 114	0.50	Not Detected	3.5	Not Detected
Chloromethane	5.0	Not Detected	10	Not Detected
Vinyl Chloride	0.50	Not Detected	1.3	Not Detected
1,3-Butadiene	0.50	Not Detected	1.1	Not Detected
Bromomethane	5.0	Not Detected	19	Not Detected
Chloroethane	2.0	Not Detected	5.3	Not Detected
Freon 11	0.50	Not Detected	2.8	Not Detected
Ethanol	2.0	Not Detected	3.8	Not Detected
Freon 113	0.50	Not Detected	3.8	Not Detected
1,1-Dichloroethene	0.50	Not Detected	2.0	Not Detected
Acetone	5.0	Not Detected	12	Not Detected
2-Propanol	2.0	Not Detected	4.9	Not Detected
Carbon Disulfide	2.0	Not Detected	6.2	Not Detected
3-Chloropropene	2.0	Not Detected	6.3	Not Detected
Methylene Chloride	5.0	Not Detected	17	Not Detected
Methyl tert-butyl ether	0.50	Not Detected	1.8	Not Detected
trans-1,2-Dichloroethene	0.50	Not Detected	2.0	Not Detected
Hexane	0.50	Not Detected	1.8	Not Detected
1,1-Dichloroethane	0.50	Not Detected	2.0	Not Detected
2-Butanone (Methyl Ethyl Ketone)	2.0	Not Detected	5.9	Not Detected
cis-1,2-Dichloroethene	0.50	Not Detected	2.0	Not Detected
Tetrahydrofuran	0.50	Not Detected	1.5	Not Detected
Chloroform	0.50	Not Detected	2.4	Not Detected
1,1,1-Trichloroethane	0.50	Not Detected	2.7	Not Detected
Cyclohexane	0.50	Not Detected	1.7	Not Detected
Carbon Tetrachloride	0.50	Not Detected	3.1	Not Detected
2,2,4-Trimethylpentane	0.50	Not Detected	2.3	Not Detected
Benzene	0.50	Not Detected	1.6	Not Detected
1,2-Dichloroethane	0.50	Not Detected	2.0	Not Detected
Heptane	0.50	Not Detected	2.0	Not Detected
Trichloroethene	0.50	Not Detected	2.7	Not Detected
1,2-Dichloropropane	0.50	Not Detected	2.3	Not Detected
1,4-Dioxane	2.0	Not Detected	7.2	Not Detected
Bromodichloromethane	0.50	Not Detected	3.4	Not Detected
cis-1,3-Dichloropropene	0.50	Not Detected	2.3	Not Detected
4-Methyl-2-pentanone	0.50	Not Detected	2.0	Not Detected
Toluene	0.50	Not Detected	1.9	Not Detected
trans-1,3-Dichloropropene	0.50	Not Detected	2.3	Not Detected
1,1,2-Trichloroethane	0.50	Not Detected	2.7	Not Detected
Tetrachloroethene	0.50	Not Detected	3.4	Not Detected
2-Hexanone	2.0	Not Detected	8.2	Not Detected



Client Sample ID: Lab Blank Lab ID#: 1210554-04A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name:	p102623	Date of Collection: NA
Dil. Factor:	1.00	Date of Analysis: 10/26/12 10:44 PM

Dil. I actor.	1.00	Date of Affaiysis. 10/20/12 10:44 FM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Dibromochloromethane	0.50	Not Detected	4.2	Not Detected
1,2-Dibromoethane (EDB)	0.50	Not Detected	3.8	Not Detected
Chlorobenzene	0.50	Not Detected	2.3	Not Detected
Ethyl Benzene	0.50	Not Detected	2.2	Not Detected
m,p-Xylene	0.50	Not Detected	2.2	Not Detected
o-Xylene	0.50	Not Detected	2.2	Not Detected
Styrene	0.50	Not Detected	2.1	Not Detected
Bromoform	0.50	Not Detected	5.2	Not Detected
Cumene	0.50	Not Detected	2.4	Not Detected
1,1,2,2-Tetrachloroethane	0.50	Not Detected	3.4	Not Detected
Propylbenzene	0.50	Not Detected	2.4	Not Detected
4-Ethyltoluene	0.50	Not Detected	2.4	Not Detected
1,3,5-Trimethylbenzene	0.50	Not Detected	2.4	Not Detected
1,2,4-Trimethylbenzene	0.50	Not Detected	2.4	Not Detected
1,3-Dichlorobenzene	0.50	Not Detected	3.0	Not Detected
1,4-Dichlorobenzene	0.50	Not Detected	3.0	Not Detected
alpha-Chlorotoluene	0.50	Not Detected	2.6	Not Detected
1,2-Dichlorobenzene	0.50	Not Detected	3.0	Not Detected
1,2,4-Trichlorobenzene	2.0	Not Detected	15	Not Detected
Hexachlorobutadiene	2.0	Not Detected	21	Not Detected

Container Type: NA - Not Applicable

Surrogates	%Recovery	меtnod Limits
Toluene-d8	99	70-130
1,2-Dichloroethane-d4	94	70-130
4-Bromofluorobenzene	96	70-130



Client Sample ID: CCV Lab ID#: 1210554-05A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: p102619 Date of Collection: NA
Dil. Factor: 1.00 Date of Analysis: 10/26/12 08:05 PM

Compound	%Recovery
Freon 12	81
Freon 114	84
Chloromethane	90
Vinyl Chloride	86
1,3-Butadiene	82
Bromomethane	88
Chloroethane	89
Freon 11	79
Ethanol	92
Freon 113	81
1,1-Dichloroethene	90
Acetone	80
2-Propanol	86
Carbon Disulfide	84
3-Chloropropene	94
Methylene Chloride	84
Methyl tert-butyl ether	79
trans-1,2-Dichloroethene	88
Hexane	91
1,1-Dichloroethane	84
2-Butanone (Methyl Ethyl Ketone)	82
cis-1,2-Dichloroethene	86
Tetrahydrofuran	85
Chloroform	81
1,1,1-Trichloroethane	84
Cyclohexane	88
Carbon Tetrachloride	82
2,2,4-Trimethylpentane	93
Benzene	83
1,2-Dichloroethane	78
Heptane	93
Trichloroethene	76
1,2-Dichloropropane	87
1,4-Dioxane	77
Bromodichloromethane	84
cis-1,3-Dichloropropene	93
4-Methyl-2-pentanone	94
Toluene	88
trans-1,3-Dichloropropene	91
1,1,2-Trichloroethane	81
Tetrachloroethene	78
2-Hexanone	81



Client Sample ID: CCV Lab ID#: 1210554-05A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: p102619 Date of Collection: NA
Dil. Factor: 1.00 Date of Analysis: 10/26/12 08:05 PM

Compound	%Recovery
Dibromochloromethane	81
1,2-Dibromoethane (EDB)	84
Chlorobenzene	84
Ethyl Benzene	91
m,p-Xylene	93
o-Xylene	95
Styrene	92
Bromoform	84
Cumene	94
1,1,2,2-Tetrachloroethane	83
Propylbenzene	89
4-Ethyltoluene	89
1,3,5-Trimethylbenzene	86
1,2,4-Trimethylbenzene	91
1,3-Dichlorobenzene	84
1,4-Dichlorobenzene	80
alpha-Chlorotoluene	105
1,2-Dichlorobenzene	83
1,2,4-Trichlorobenzene	76
Hexachlorobutadiene	83

Container Type: NA - Not Applicable

		Method Limits	
Surrogates	%Recovery		
Toluene-d8	104	70-130	
1,2-Dichloroethane-d4	94	70-130	
4-Bromofluorobenzene	103	70-130	



Client Sample ID: LCS Lab ID#: 1210554-06A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: p102621 Date of Collection: NA
Dil. Factor: 1.00 Date of Analysis: 10/26/12 08:57 PM

Compound	%Recovery
Freon 12	85
Freon 114	88
Chloromethane	89
Vinyl Chloride	92
1,3-Butadiene	86
Bromomethane	91
Chloroethane	93
Freon 11	82
Ethanol	60 Q
Freon 113	86
1,1-Dichloroethene	101
Acetone	84
2-Propanol	90
Carbon Disulfide	111
3-Chloropropene	117
Methylene Chloride	87
Methyl tert-butyl ether	74
trans-1,2-Dichloroethene	103
Hexane	100
1,1-Dichloroethane	88
2-Butanone (Methyl Ethyl Ketone)	84
cis-1,2-Dichloroethene	92
Tetrahydrofuran	86
Chloroform	86
1,1,1-Trichloroethane	90
Cyclohexane	94
Carbon Tetrachloride	87
2,2,4-Trimethylpentane	97
Benzene	87
1,2-Dichloroethane	80
Heptane	96
Trichloroethene	79
1,2-Dichloropropane	90
1,4-Dioxane	77
Bromodichloromethane	87
cis-1,3-Dichloropropene	98
4-Methyl-2-pentanone	94
Toluene	90
trans-1,3-Dichloropropene	97
1,1,2-Trichloroethane	85
Tetrachloroethene	81
2-Hexanone	81



Client Sample ID: LCS Lab ID#: 1210554-06A

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: p102621 Date of Collection: NA
Dil. Factor: 1.00 Date of Analysis: 10/26/12 08:57 PM

Compound	%Recovery
Dibromochloromethane	84
1,2-Dibromoethane (EDB)	90
Chlorobenzene	89
Ethyl Benzene	95
m,p-Xylene	99
o-Xylene	99
Styrene	96
Bromoform	86
Cumene	98
1,1,2,2-Tetrachloroethane	89
Propylbenzene	95
4-Ethyltoluene	92
1,3,5-Trimethylbenzene	89
1,2,4-Trimethylbenzene	94
1,3-Dichlorobenzene	89
1,4-Dichlorobenzene	85
alpha-Chlorotoluene	109
1,2-Dichlorobenzene	89
1,2,4-Trichlorobenzene	83
Hexachlorobutadiene	89

Q = Exceeds Quality Control limits.

Container Type: NA - Not Applicable

	Method	
%Recovery	Limits	
103	70-130	
94	70-130	
102	70-130	
	103 94	



Client Sample ID: LCSD Lab ID#: 1210554-06AA

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: p102622 Date of Collection: NA
Dil. Factor: 1.00 Date of Analysis: 10/26/12 10:08 PM

Compound	%Recovery
Freon 12	87
Freon 114	91
Chloromethane	97
Vinyl Chloride	96
1,3-Butadiene	87
Bromomethane	94
Chloroethane	95
Freon 11	84
Ethanol	61 Q
Freon 113	88
1,1-Dichloroethene	101
Acetone	85
2-Propanol	92
Carbon Disulfide	115
3-Chloropropene	114
Methylene Chloride	87
Methyl tert-butyl ether	76
trans-1,2-Dichloroethene	106
Hexane	103
1,1-Dichloroethane	91
2-Butanone (Methyl Ethyl Ketone)	87
cis-1,2-Dichloroethene	93
Tetrahydrofuran	87
Chloroform	88
1,1,1-Trichloroethane	92
Cyclohexane	96
Carbon Tetrachloride	89
2,2,4-Trimethylpentane	98
Benzene	88
1,2-Dichloroethane	82
Heptane	98
Trichloroethene	80
1,2-Dichloropropane	92
1,4-Dioxane	80
Bromodichloromethane	90
cis-1,3-Dichloropropene	99
4-Methyl-2-pentanone	92
Toluene	91
trans-1,3-Dichloropropene	95
1,1,2-Trichloroethane	84
Tetrachloroethene	80
2-Hexanone	80



Client Sample ID: LCSD Lab ID#: 1210554-06AA

EPA METHOD TO-15 GC/MS FULL SCAN

File Name: p102622 Date of Collection: NA
Dil. Factor: 1.00 Date of Analysis: 10/26/12 10:08 PM

Compound	%Recovery
Dibromochloromethane	84
1,2-Dibromoethane (EDB)	88
Chlorobenzene	88
Ethyl Benzene	94
m,p-Xylene	98
o-Xylene	99
Styrene	96
Bromoform	86
Cumene	97
1,1,2,2-Tetrachloroethane	88
Propylbenzene	94
4-Ethyltoluene	88
1,3,5-Trimethylbenzene	88
1,2,4-Trimethylbenzene	92
1,3-Dichlorobenzene	87
1,4-Dichlorobenzene	83
alpha-Chlorotoluene	109
1,2-Dichlorobenzene	86
1,2,4-Trichlorobenzene	80
Hexachlorobutadiene	85

Q = Exceeds Quality Control limits.

Container Type: NA - Not Applicable

		Method
Surrogates	%Recovery	Limits
Toluene-d8	103	70-130
1,2-Dichloroethane-d4	94	70-130
4-Bromofluorobenzene	101	70-130

CHAIN-OF-CUSTODY RECORD

Sample Transportation Notice
Relinquishing signature on this document indicates that sample is being shipped in compliance with all applicable local, State, Federal, national, and international laws, regulations and ordinances of any kind. Air Toxics Limited assumes no liability with respect to the collection, handling or shipping of these samples. Relinquishing signature also indicates agreement to hold harmless, defend, and indemnify Air Toxics Limited against any claim, demand, or action, of any kind, related to the collection, handling, or shipping of samples. D.O.T. Hotline (800) 467-4922

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DATA USABILITY SUMMARY REPORT

LUDLOW FORMER MGP SITE

SUPPLEMENTAL REMEDIAL INVESTIGATION

Prepared For:



CONSOLIDATED EDISON COMPANY OF NEW YORK, INC.

31-01 20th Avenue Long Island City, NY 11105

Prepared By:

PARSONS

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FEBRUARY 2013

TABLE OF CONTENTS

<u>Pa</u>	<u>ge</u>
SECTION 1 DATA USABILITY SUMMARY1	-1
1.2 SAMPLING AND CHAIN-OF-CUSTODY	1
1.3 LABORATORY ANALYTICAL METHODS1	-1
1.3.1 Volatile Organic Analysis	-2
1.3.2 Semivolatile Organic Analysis	
1.3.3 Inorganics Analysis	
SECTION 2 DATA VALIDATION REPORT2	2-1
2.1 SOIL	2-1
2.1.1 Volatiles	
2.1.2 Semivolatiles	2-4
2.1.3 Inorganics	2-5
2.2 GROUNDWATER	2-7
2.2.1 Volatiles	2-7
2.2.2 Semivolatiles	2-9
2.2.3 Inorganics	10
LIST OF ATTACHMENTS	
ATTACHMENT A VALIDATED LABORATORY DATA	
ATTACHMENT A-1 VALIDATED LABORATORY DATA FOR SOIL SA	MPLES
ATTACHMENT A-2 VALIDATED LABORATORY DATA FOR	
GROUNDWATER SAMPLES	

SECTION 1

DATA USABILITY SUMMARY

Soil and groundwater samples were collected from the Consolidated Edison Ludlow Street Site from October 25, 2012 through December 21, 2012. Analytical results from these samples were validated and reviewed by Parsons for usability with respect to the following requirements:

- Work Plan.
- NYSDEC Analytical Services Protocol (ASP), and
- USEPA Region II Standard Operating Procedures (SOPs) for organic and inorganic data review.

The analytical laboratory for this project was Chemtech. This laboratory is certified to perform project analyses through the New York State Department of Health (NYSDOH) Environmental Laboratory Approval Program (ELAP).

1.1 LABORATORY DATA PACKAGES

The laboratory data package turnaround time, defined as the time from sample receipt by the laboratory to receipt of the analytical data packages by Parsons, was 16-27 days for the project samples.

The data packages received from Chemtech were paginated, complete, and overall were of good quality. Comments on specific quality control (QC) and other requirements are discussed in detail in the attached data validation report which is summarized by media in Section 2.

1.2 SAMPLING AND CHAIN-OF-CUSTODY

The samples were collected, properly preserved, shipped under a chain-of-custody (COC) record, and received at Chemtech within one to seven days of sampling. All samples were received intact and in good condition at Chemtech.

1.3 LABORATORY ANALYTICAL METHODS

The soil and groundwater samples that were collected from the site were analyzed for volatile organic compounds (VOCs), semivolatile organic compounds (SVOCs), metals, and cyanide. Summaries of issues concerning these laboratory analyses are presented in Subsections 1.3.1 through 1.3.3. The data qualifications resulting from the data validation review and statements on the laboratory analytical precision, accuracy, representativeness, completeness, and comparability (PARCC) are discussed for each analytical method in Section 2. The laboratory data were reviewed and may be qualified with the following validation flags:

"U" - not detected at the value given,

"UJ" - estimated and not detected at the value given,

"J" - estimated at the value given,

"N" - presumptive evidence at the value given, and

"R" - unusable value.

The validated laboratory data were tabulated and are presented in Attachment A.

1.3.1 Volatile Organic Analysis

Soil and groundwater samples were analyzed for VOCs using the USEPA SW-846 8260C analytical method. Certain reported results for the VOC samples were qualified as estimated based upon instrument calibrations, laboratory control sample recoveries, internal standard responses, and field duplicate precision. Certain reported VOC analytical results were considered unusable and qualified "R" based upon poor instrument calibration linearity and internal standard responses. The reported VOC analytical results were 98.8% and 100% complete (i.e., usable) for the soil and groundwater data, respectively. PARCC requirements were met overall.

1.3.2 Semivolatile Organic Analysis

Soil and groundwater samples were analyzed for SVOCs using the USEPA SW-846 8270D analytical method. Certain reported results for the SVOC samples were qualified as estimated based upon matrix spike/matrix spike duplicate precision and accuracy, laboratory control sample recoveries, instrument calibrations, and internal standard responses. The reported SVOC analytical results were 100% complete (i.e., usable) for the soil and groundwater data. PARCC requirements were met.

1.3.3 Inorganics Analysis

Soil and groundwater samples were analyzed for metals and cyanide using the USEPA SW-846 6010B/7470A/7471A/9012B analytical methods. Certain reported results for the inorganics samples were qualified as estimated based upon matrix spike recoveries, serial dilutions, laboratory duplicate precision, and field duplicate precision. The reported inorganic analytical results were considered 100% complete (i.e., usable) for the soil and groundwater data. PARCC requirements were met.

SECTION 2

DATA VALIDATION REPORT

2.1 SOIL

Data review has been completed for data packages generated by Chemtech containing soil samples collected from the site. All of these samples were properly preserved, shipped under a COC record, and received intact by the analytical laboratory. The analytical results were presented by the laboratory in three sample delivery groups (SDGs): D4710, D4751, and D5300. Data validation was performed for all samples in accordance with the most current editions of the USEPA Region II SOPs for organic and inorganic data review. This data validation and usability report is presented by analysis type and the validated laboratory data are presented in Attachment A-1.

2.1.1 Volatiles

The following items were reviewed for compliancy in the volatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- Matrix spike/matrix spike duplicate (MS/MSD) precision and accuracy
- Laboratory control sample (LCS) recoveries
- Laboratory method blank and field equipment blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times.
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, blank contamination, initial and continuing calibrations, internal standard responses, and field duplicate precision as discussed below.

MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC acceptance limits for designated

project spiked samples with the exception of many precision outliers during the spiked analyses of sample MW-7(48-50). Validation qualification of the parent sample was not required since the MS/MSD accuracy measurements were within criteria.

LCS Recoveries

All LCS recoveries associated with project samples were considered acceptable and within QC limits with the exception of the high LCS recoveries for methylene chloride (140%R, 175%R; QC limit 73-134%R) associated with all samples in SDG D4710 except SB-10(45-47); the high LCS recovery for acetone (160%R; QC limit 57-148%R) associated with sample SB-12(12-14); the low LCS recovery for bromochloromethane (75%R; QC limit 76-123%R) associated with all samples in SDG D4751 except SB-12(12-14); the high LCS recoveries for 1,2,4-trichlorobenzene (135%R; QC limit 65-130%R) and 1,2,3-trichlorobenzene (140%R; QC limit 60-135%R) associated with samples SB-16(45-47), SB-14(45-47), SB-14A(45-47), and SB-17(33-35); the low LCS recovery for 1,2,3-trichlorobenzene (70%R; QC limit 79-123%R) and the high LCS recovery for bromomethane (140%R; QC limit 64-136%R) associated with samples SB-14(38-40) and SB-17(45-47). Therefore, positive results for those compounds where the LCS recovery exceeded the QC limit were considered estimated, possibly biased high, and qualified "J" for the affected samples. Results for those compounds where the LCS recovery fell below the QC limit were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

Blank Contamination

The laboratory method blanks associated with samples in SDGs D4710 and D4751 contained methylene chloride at concentrations ranging from 4.6 to 9.5 μ g/kg. Therefore, all methylene chloride results less than validation action concentrations were considered not detected and qualified "U" for the affected samples.

Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of methylene chloride (24.8%RSD) and 1,2,3-trichlorobenzene (21.3%RSD) in the initial calibration associated with samples SB-14(38-40) and SB-17(45-47); carbon tetrachloride (36.9%RSD) in the initial calibration associated with sample FB122112; 2-hexanone (21.2%RSD), dibromochloromethane (26.5%RSD), and bromoform (34.5%RSD) in the initial calibration associated with samples SB-16(30-33) and –DL; and 1,4-dioxane (31.3%RSD) and 1,2-dibromo-3-chloropropane (RRF=0.048) in the initial calibration associated with samples SB-16(45-47), SB-14(45-47), SB-14A(45-47), and SB-17(33-35). The results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples. However, the results for those compounds where the RRF was outside the criteria which were nondetects were considered unusable and qualified "R" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within $\pm 20\%$ with the exception of methylene chloride (32.39%D, 38.81%D, 77.31%D) in the continuing calibrations associated with samples in SDG

D4710; methylene chloride (38.36%D) and 1,2-dibromo-3-chloropropane (-20.25%D) in the continuing calibration associated with all samples in SDG D4751 except SB-11(29.5-30), SB-12(12-14), "RE" samples, and "DL" samples; acetone (20.36%D) and methylene chloride (21.19%D) in the continuing calibration associated with samples SB-11(29.5-30), MW-8(9.5-10)RE, and TP-6FLOORRE; chloroethane (30.46%D), acetone (45.68%D), methyl acetate (38.82%D), and 2-hexanone (21.05%D) in the continuing calibration associated with sample SB-12(12-14); bromomethane (49.57%D) and methylene chloride (-20.25%D) in the continuing calibration associated with samples SB-14(38-40) and SB-17(45-47); 2-butanone (20.99%D) and 1,4-dioxane (50%D) in the continuing calibration associated with samples SB-16(30-33) and -DL; dichlorodifluoromethane (-20.07%D), dibromochloromethane (21.33%D), and 1,2,4trichlorobenzene (26.25%D) in the continuing calibration associated with samples SB-16(45-47), SB-14(45-47), SB-14A(45-47), and SB-17(33-35); and bromomethane (68.79%D), carbon tetrachloride (65.34%D), and 1,4-dioxane (-33.33%D) in the continuing calibration associated with FB122112. Therefore, the sample results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

Internal Standard Responses

All internal standard (IS) responses and retention times were within specified QC ranges based on associated calibration standards (i.e., sample's area count within -50% to +100% and retention times within ±0.5 minutes of the standard) with the exception of the low response for the IS pentafluorobenzene in sample SB-10(30-32)DL; low response for the IS chlorobenzene-d5 in samples SB-9(5-8), TP-6FLOOR, and -RE; low response for the IS 1,4-difluorobenzene-d4 in samples SB-9(5-8), -RE, MW-17(36-38), MW-7(36-38), MW-8(9.5-10), and -RE; and the extremely low response for the IS 1,4-difluorobenzene-d4 in samples SB-10(30-32), TP-6FLOOR, and -RE. Therefore, results associated with these ISs were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples. However, nondetected results associated with those ISs that experienced extremely low responses were considered unusable and qualified "R" for the affected samples.

Field Duplicate Precision

All field duplicate precision results were considered acceptable with the exception of the precision results for cyclohexane (153%RPD) and methylcyclohexane (128%RPD) for the field duplicate pair MW-7(36-38) and MW-17(36-38). Therefore, the results for these compounds were considered estimated and qualified "J" for these samples.

Usability

All volatile soil sample results were considered usable following data validation with the exception of certain nondetected results based upon poor instrument calibration linearity and internal standard responses.

Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The volatile soil data presented

by Chemtech were 98.8% complete (i.e., usable). The validated volatile laboratory data are tabulated and presented in Attachment A-1.

2.1.2 Semivolatiles

The following items were reviewed for compliancy in the semivolatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- MS/MSD precision and accuracy
- LCS recoveries
- Laboratory method blank and field equipment blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, and initial and continuing calibrations as discussed below.

MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated spiked project samples with the exception of the high MS/MSD accuracy results for hexachlorocyclopentadiene (130%R/130%R; QC limit 10-122%R) during the spiked analyses of sample MW-7(48-50). Validation qualification of the parent sample was not required since this compound was not detected.

LCS Recoveries

All LCS recoveries were considered acceptable and within QC limits with the exception of the high LCS recoveries for 2-nitrophenol (106%R; QC limit 52-105%R), hexachlorocyclopentadiene (161%R; QC limit 38-122%R), 2,4,6-trichlorophenol (106%R; QC limit 56-103%R), acenaphthene (106%R; QC limit 57-102%R), benzo(b)fluoranthene (106%R; QC limit 56-103%R), benzo(a)pyrene (112%R; QC limit 57-103%R), and benzo(g,h,i)perylene (106%R; QC limit 56-105%R) associated with samples in SDG D4710; the LCS recoveries for 4-chloroaniline (16%R; QC limit 25-115%R) and hexachlorocyclopentadiene (142%R;43-

112%R) associated with samples **SDG** D4751; the LCS recoveries for in hexachlorocyclopentadiene (27%R; QC limit 38-122%R), 2,4-dinitrophenol (15%R; QC limit 32-114%R), and 4,6-dinitro-2-methylphenol (18%R; QC limit 44-119%R) associated with soil samples in SDG D5300; and the LCS recovery for hexachlorocyclopentadiene (130%R; QC limit 42-121%R) associated with sample FB122112. Therefore, results for those compounds where LCS recoveries fell below the QC limit were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples. Positive results for those compounds where the LCS recovery exceeded the QC limit were considered estimated, possibly biased high, and qualified "J" for the affected samples.

Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of 2,4-dinitrophenol (62.5%RSD) and 4,6-dinitro-2-methylphenol (39.2%RSD) in the initial calibration associated with all samples in SDG D4751 except SB-11(29.5-30), MW-8(24.5-25), and MW-9(35-35.5); 2,4-dinitrophenol (24.5%RSD) in the initial calibration associated with samples SB-11(29.5-30), MW-8(24.5-25), and MW-9(35-35.5); and 2,4-dinitrophenol (38.1%RSD) and 4,6-dinitro-2-methylphenol (23.6%RSD) in the initial calibration associated with samples in SDG D5300. The results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within $\pm 20\%$ with the exception of 2,4-dinitrophenol (86.5%D) and 4,6-dinitro-2-methylphenol (21.7%D) in the continuing calibration associated with sample SB-12(36-38). Therefore, results for these compounds which were nondetects were considered estimated and qualified "UJ" for this sample.

Usability

All semivolatile soil sample results were considered usable following data validation.

Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The semivolatile soil data presented by Chemtech were 100% complete (i.e., usable). The validated semivolatile laboratory data are tabulated and presented in Attachment A-1.

2.1.3 Inorganics

The following items were reviewed for compliancy in the inorganics analysis:

- Custody documentation
- Holding times
- Initial and continuing calibration verifications

- Initial and continuing calibration blank, and laboratory preparation blank, and field equipment blank contamination
- Inductively coupled plasma (ICP) interference check sample (ICS)
- Matrix spike (MS) recoveries
- Laboratory duplicate precision
- Field duplicate precision
- Laboratory control sample (LCS) recoveries
- ICP serial dilutions
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of blank contamination, matrix spike recoveries, and serial dilutions as discussed below.

Blank Contamination

The field equipment blank FB122112 associated with samples in SDG D5300 contained chromium, iron, and potassium at concentrations of 21.3, 551, and 1190 μ g/L, respectively. Therefore, all results for these analytes less than the validation action concentrations were considered not detected and qualified "U" for the affected samples.

Matrix Spike Recoveries

All the MS recoveries for designated spiked project samples were within the 75-125%R QC limit with sample concentrations less than four times the spiking concentration with the exception of the low MS recoveries for antimony (52.1%R, 51.2%R, 74.1%R, 74.3%R) associated with all soil samples in SDGs D4751 and D5300. Therefore, results for antimony were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples since MS recoveries fell below the QC limit.

ICP Serial Dilutions

All serial dilution results for designated project samples were considered acceptable with a percent difference (%D) less than 10% for all ICP analytes with the exception of chromium (15.8%D) and copper (12.5%D) associated with samples in SDG D4710; aluminum (14.8%D), arsenic (17%D), barium (12%D), calcium (14.3%D), chromium (18.9%D), copper (23.4%D), magnesium (16.2%D), manganese (30.3%D), potassium (12.8%D), sodium (14%D), and vanadium (19.4%D) associated with samples in SDG D4751; and aluminum (13.9%D), barium (25.1%D), calcium (28.9%D), chromium (26.6%D), iron (13.9%D), magnesium (23.8%D), manganese (29.1%D), potassium (25.2%D), and zinc (15.4%D) associated with soil samples in

SDG D5300. Therefore, positive results for these analytes were considered estimated and qualified "J" for the affected samples.

<u>Usability</u>

All inorganics soil sample results were considered usable following data validation.

Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The inorganics soil data presented by Chemtech were 100% complete (i.e., usable). The validated soil inorganics laboratory data are tabulated and presented in Attachment A-1.

2.2 GROUNDWATER

Data review has been completed for data packages generated by Chemtech containing groundwater samples collected from the site. All of these samples were properly preserved, shipped under a COC record, and received intact by the analytical laboratory. The analytical results were presented by the laboratory in one sample delivery groups (SDGs): D4947. Data validation was performed for all samples in accordance with the most current editions of the USEPA Region II SOPs for organic and inorganic data review. This data validation and usability report is presented by analysis type and the validated laboratory data are presented in Attachment A-2.

2.2.1 Volatiles

The following items were reviewed for compliancy in the volatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- Matrix spike/matrix spike duplicate (MS/MSD) precision and accuracy
- Laboratory control sample (LCS) recoveries
- Laboratory method blank and field equipment/trip blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, blank contamination, and initial and continuing calibrations as discussed below.

MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated project spiked samples with the exception of the high accuracy results for carbon tetrachloride (150%R/140%R; QC limit 60-140%R) during the spiked analyses of sample MW-9. Since this compound was not detected in the unspiked parent sample, validation qualification was not required.

Blank Contamination

The field equipment blank FB112712 associated with groundwater samples contained carbon disulfide and methylene chloride at concentrations of 2 and 4.3 μ g/L, respectively. Since these compounds were not detected in the project samples, validation qualification was not required.

Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of acetone (20.4%RSD), chloromethane (21.6%RSD), bromomethane (21.4%RSD), carbon tetrachloride (67.5%RSD), dibromochloromethane (21.9%RSD), and bromoform (27.8%RSD) in the initial calibration associated with all samples except MW-4, -7DL, and -3DL. The results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within ±20% with the exception of methyl acetate (20.78%D), carbon tetrachloride (45.25%D), bromodichloromethane (21.58%D), 4-methyl-2-pentanone (20.34%D), trans-1,3-dichloropropene (23.81%D), dibromochloromethane (30.94%D), bromoform (43.33%D), and 1,4-dioxane (33.33%D) in the continuing calibration associated with all groundwater samples except MW-4, -3DL, and -7DL. Therefore, the sample results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

Usability

All volatile groundwater sample results were considered usable following data validation.

Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The volatile groundwater data presented by Chemtech were 100% complete (i.e., usable). The validated volatile laboratory data are tabulated and presented in Attachment A-2.

2.2.2 Semivolatiles

The following items were reviewed for compliancy in the semivolatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- MS/MSD precision and accuracy
- LCS recoveries
- Laboratory method blank and field equipment blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, initial and continuing calibrations, and internal standard responses as discussed below.

MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated spiked project samples with the exception of the low MS/MSD accuracy results for 2,3,4,6-tetrachlorophenol (85%R/86%R; QC limit 91-111%R) during the spiked analyses of sample MW-9. Therefore, the nondetected result for this compound was considered estimated, possibly biased low, and qualified "UJ" for the parent sample.

LCS Recoveries

All LCS recoveries were considered acceptable and within QC limits with the exception of the high LCS recoveries for 2-methylphenol (100%R; QC limit 32-94%R), 3+4-methylphenols (99%R; QC limit 24-91%R), hexachlorocyclopentadiene (150%R; QC limit 42-121%R), and 1,2,4,5-tetrachlorobenzene (106%R; QC limit 60-105%R) associated with all groundwater samples. Validation qualification of the groundwater samples was not required since these compounds were not detected.

Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of 2,4-dinitrophenol (43.3%RSD, 34.8%RSD) and 4,6-dinitro-2-methylphenol (25.1%RSD, 23.4%RSD) in the initial calibrations associated with all samples. The results for these compounds which were nondetects were considered estimated and qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within ±20% with the exception of 2,4-dinitrophenol (34.4%D) and 4,6-dinitro-2-methylphenol (23.6%D) in the continuing calibration associated with samples MW-7, 18, -8, -1, and -9; and 2,4-dinitrophenol (21.6%D) in the continuing calibration associated with sample FB112712. Therefore, the sample results for these compounds which were nondetects were considered estimated and qualified "UJ" for the affected samples.

Internal Standard Responses

All internal standard (IS) responses and retention times were within specified QC ranges based on associated calibration standards (i.e., sample's area count within -50% to +100% and retention times within ± 0.5 minutes of the standard) with the exception of the low responses for the ISs acenaphthene-d10 and phenanthrene-d10 in sample MW-3. Therefore, results associated with these ISs were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected sample.

Usability

All semivolatile groundwater sample results were considered usable following data validation.

Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The semivolatile groundwater data presented by Chemtech were 100% complete (i.e., usable). The validated semivolatile laboratory data are tabulated and presented in Attachment A-2.

2.2.3 Inorganics

The following items were reviewed for compliancy in the inorganics analysis:

- Custody documentation
- Holding times
- Initial and continuing calibration verifications
- Initial and continuing calibration blank, and laboratory preparation blank, and field equipment blank contamination
- Inductively coupled plasma (ICP) interference check sample (ICS)

- Matrix spike (MS) recoveries
- Laboratory duplicate precision
- Field duplicate precision
- Laboratory control sample (LCS) recoveries
- ICP serial dilutions
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of blank contamination, laboratory duplicate precision, field duplicate precision, and serial dilutions as discussed below.

Blank Contamination

The field equipment blank FB112712 associated with the groundwater samples contained aluminum, chromium, iron, nickel, potassium, sodium, and zinc at concentrations of 61.4, 25.7, 161, 10.5, 763, 17,600, and 14 μ g/L, respectively. Therefore, all results for these analytes less than the validation action concentrations were considered not detected and qualified "U" for the affected samples.

Laboratory Duplicate Precision

All laboratory duplicate precision results were considered acceptable with precision results less than 20%RPD for all analytes with the exception of aluminum (33.5%RPD), antimony (20.7%RPD), barium (42.8%RPD), beryllium (23.6%RPD), calcium (45.7%RPD), chromium (42.1%RPD), copper (24.1%RPD), iron (43.2%RPD), magnesium (46.1%RPD), manganese (40.5%RPD), potassium (29%RPD), silver (23.6%RPD), sodium (28.7%RPD), thallium (26.6%RPD), vanadium (36.3%RPD), and zinc (82.8%RPD) associated with all groundwater samples. Therefore, the results for these analytes were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

Field Duplicate Precision

All field duplicate precision results were considered acceptable for the field duplicate pair MW-8 and MW-18 with the exception of the precision for aluminum (133%RPD), arsenic (104%RPD), cobalt (99%RPD), copper (96%RPD), iron (133%RPD), lead (111%RPD), nickel (52%RPD), vanadium (117%RPD), and zinc (99%RPD). Therefore, the results for these analytes were considered estimated and qualified "J" for the affected parent sample and its field duplicate.

ICP Serial Dilutions

All serial dilution results for designated project samples were considered acceptable with a percent difference (%D) less than 10% for all ICP analytes with the exception of manganese

(11.7%D) and sodium (74.6%D) associated with groundwater samples. Therefore, positive results for these analytes were considered estimated and qualified "J" for the affected samples.

Usability

All inorganics groundwater sample results were considered usable following data validation.

Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The inorganics groundwater data presented by Chemtech were 100% complete (i.e., usable). The validated groundwater inorganics laboratory data are tabulated and presented in Attachment A-2.

ATTACHMENT A VALIDATED LABORATORY DATA

ATTACHMENT A-1 VALIDATED LABORATORY DATA FOR SOIL SAMPLES

				Dup of MW-7(36-38)		
Con Ed - Lud	ow Site	Location ID:	MW- 7	MW- 7	MW- 7	MW-8
Validated Soil	Analytical Data	Sample ID:	MW-7(36-38)	MW-17(36-38)	MW-7(48-50)	MW-8(9.5-10)
SDGs: D4710	, D1751, D5300	Lab Sample Id:	D4710-03	D4710-09	D4710-04	D4751-03
		Depth:	36 - 38 ft	36 - 38 ft	48 - 50 ft	9.5 - 10 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4710	D4710	D4710	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	10/26/2012 9:10	10/26/2012 9:15	10/26/2012 9:20	11/6/2012 14:15
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
71-55-6	VOLATILES 1,1,1-TRICHLOROETHANE	na/ka	5.4 U	1.1 U	0.96 U	1.1 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg ug/kg	2.8 UJ	0.55 UJ	0.5 U	0.56 UJ
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	8.2 U	1.6 U	1.5 U	1.6 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	5.6 U	1.1 U	0.98 U	1.1 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	5.8 U	1.1 U	1 U	1.1 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	9.1 U	1.8 U	1.6 U	1.8 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	3.1 UJ	0.6 UJ	0.55 U	0.61 UJ
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	4.3 UJ	0.84 UJ	0.76 U	0.86 UJ
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	5.4 UJ	1 UJ	0.95 U	1.1 UJ
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/kg	4 U	0.77 U	0.7 U	0.78 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	3.8 UJ	0.75 UJ	0.68 U	0.76 UJ
107-06-2	1,2-DICHLOROETHANE	ug/kg	4 U	0.77 U	0.7 U	0.78 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	1.6 U	0.31 U	0.28 U	0.32 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	2.3 UJ	0.44 UJ	0.4 U	0.45 UJ
106-46-7	1,4-DICHLOROBENZENE	ug/kg	2.5 UJ	0.49 UJ	0.45 U	0.5 UJ
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	310 U	60 U	55 U	60 U
591-78-6	2-HEXANONE	ug/kg	24 U	4.7 U	4.3 U	4.8 U
67-64-1	ACETONE	ug/kg	19 U	3.6 U	3.3 U	51
71-43-2	BENZENE BROMOGUL OROMETHANE	ug/kg	7100 J	5300	1.3 J	2.3 J
74-97-5	BROMOCHLOROMETHANE	ug/kg	4.9 U	0.95 U	0.86 U	0.97 U
75-27-4 75-25-2	BROMODICHLOROMETHANE BROMOFORM	ug/kg	3.8 U 4.6 U	0.75 U 0.89 U	0.68 U 0.81 U	0.76 U 0.9 U
74-83-9	BROMOMETHANE	ug/kg ug/kg	4.6 U	2.9 U	2.7 U	0.9 U
75-15-0	CARBON DISULFIDE	ug/kg ug/kg	6.5 U	1.3 U	1.3 J	4.9 J
56-23-5	CARBON TETRACHLORIDE	ug/kg	6.1 U	1.2 U	1.1 U	1.2 U
108-90-7	CHLOROBENZENE	ug/kg	3.1 U	0.6 U	0.55 U	0.61 U
75-00-3	CHLOROETHANE	ug/kg	8.6 U	1.7 U	1.5 U	1.7 U
67-66-3	CHLOROFORM	ug/kg	4.6 U	0.89 U	0.81 U	0.9 U
74-87-3	CHLOROMETHANE	ug/kg	5.3 U	1 U	0.94 U	1.1 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	5.5 U	1.1 U	0.97 U	1.1 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	4.4 U	0.87 U	0.79 U	0.88 U
110-82-7	CYCLOHEXANE	ug/kg	62 J	8.2 J	1.1 U	1.2 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	3.3 U	0.65 U	0.59 U	0.66 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	4 U	0.78 U	0.71 U	3.4 J
100-41-4	ETHYLBENZENE	ug/kg	30000 J	48000	1.7 J	0.76 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	6200 J	3600	0.52 U	0.59 UJ
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	34000 J	54000	2 J	0.88 U
79-20-9	METHYL ACETATE	ug/kg	9.3 U	1.8 U	1.6 U	1.8 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/kg	19 U	3.7 U	3.4 U	3.8 U
108-10-1 108-87-2	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/kg	18 U	3.5 U	3.2 U	3.6 U
75-09-2	METHYLCYCLOHEXANE METHYLENE CHLORIDE	ug/kg	210 J 42 UJ	46 J 20 UJ	1.2 U 11 UJ	1.3 U 6.1 UJ
75-09-2 95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg ug/kg	18000 J	28000	0.74 U	0.83 U
100-42-5	STYRENE	ug/kg ug/kg	890	960 J	0.74 U 0.49 U	0.55 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg ug/kg	5.9 U	1.2 U	1 U	1.2 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	6.2 U	1.2 U	1.1 U	1.2 U
108-88-3	TOLUENE	ug/kg	15000 J	18000	1.3 J	0.78 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	4.3 U	0.83 U	0.75 U	0.84 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	4.9 U	0.95 U	0.86 U	0.97 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	5.3 U	1 U	0.94 U	1.1 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg	8.1 U	1.6 U	1.4 U	1.6 U
75-01-4	VINYL CHLORIDE	ug/kg	7.6 U	1.5 U	1.3 U	1.5 U

		Dup of MW-7(36-38)								
Con Ed - Ludlo		Location ID:	MW- 7	MW- 7	MW- 7	MW-8				
Validated Soil A		Sample ID:	MW-7(36-38)	MW-17(36-38)	MW-7(48-50)	MW-8(9.5-10)				
SDGs: D4710, I	D1751, D5300	Lab Sample Id:	D4710-03	D4710-09	D4710-04	D4751-03				
		Depth:	36 - 38 ft	36 - 38 ft	48 - 50 ft	9.5 - 10 ft				
		Source: SDG:	CTECH D4710	CTECH D4710	CTECH D4710	CTECH D4751				
		Matrix:	SOIL	SOIL	SOIL	SOIL				
		Sampled:	10/26/2012 9:10	10/26/2012 9:15	10/26/2012 9:20	11/6/2012 14:15				
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013				
CAS NO.	COMPOUND	UNITS:								
	SEMIVOLATILES	ļ								
	1,2,4,5-TETRACHLOROBENZENE	ug/kg	16 U	16 U	14 U	16 U				
	2,3,4,6-TETRACHLOROPHENOL 2,4,5-TRICHLOROPHENOL	ug/kg	16 U 29 U	16 U 28 U	14 U 26 U	16 U 28 U				
	2,4,6-TRICHLOROPHENOL	ug/kg ug/kg	13 U	12 U	11 U	12 U				
	2,4-DICHLOROPHENOL	ug/kg	16 U	15 U	14 U	15 U				
	2,4-DIMETHYLPHENOL	ug/kg	23 U	23 U	21 U	23 U				
	2,4-DINITROPHENOL	ug/kg	42 U	41 U	37 U	41 UJ				
	2,4-DINITROTOLUENE	ug/kg	12 U	12 U	11 U	12 U				
	2,6-DINITROTOLUENE	ug/kg	17 U	16 U	15 U	17 U				
	2-CHLORONAPHTHALENE 2-CHLOROPHENOL	ug/kg ug/kg	9.4 U 22 U	9.1 U 21 U	8.3 U 19 U	9.2 U 21 U				
	2-METHYLNAPHTHALENE	ug/kg ug/kg	110000	130000	9.2 U	10 U				
	2-METHYLPHENOL (O-CRESOL)	ug/kg	22 U	22 U	20 U	22 U				
	2-NITROANILINE	ug/kg	18 U	18 U	16 U	18 U				
	2-NITROPHENOL	ug/kg	20 U	19 U	18 U	20 U				
	3,3'-DICHLOROBENZIDINE	ug/kg	26 U	26 U	23 U	26 U				
	3+4-Methylphenols	ug/kg	21 U	21 U	19 U	21 U				
	3-NITROANILINE	ug/kg	26 U	26 U 23 U	23 U	26 U 23 UJ				
	4,6-DINITRO-2-METHYLPHENOL 4-BROMOPHENYL PHENYL ETHER	ug/kg ug/kg	24 U 8 U	7.8 U	21 U 7.1 U	7.9 U				
	4-CHLORO-3-METHYLPHENOL	ug/kg ug/kg	18 U	18 U	16 U	18 U				
	4-CHLOROANILINE	ug/kg	29 U	28 U	26 U	29 UJ				
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	22 U	22 U	20 U	22 U				
	4-NITROANILINE	ug/kg	53 U	52 U	48 U	53 U				
	4-NITROPHENOL	ug/kg	76 U	74 U	68 U	75 U				
	ACENAPHTHENE	ug/kg	15000 J 12000	21000 J 16000	10 U 9.2 U	11 U 10 U				
	ACENAPHTHYLENE ACETOPHENONE	ug/kg ug/kg	12000 13 U	16000 12 U	9.2 U 11 U	10 U				
	ANTHRACENE	ug/kg ug/kg	13000	20000	7.5 U	8.3 U				
	ATRAZINE	ug/kg	22 U	21 U	19 U	21 U				
100-52-7	BENZALDEHYDE	ug/kg	21 U	21 U	19 U	21 U				
	BENZO(A)ANTHRACENE	ug/kg	8300	13000	17 U	260 J				
	BENZO(A)PYRENE	ug/kg	5700 J	9000 J	7.9 U	260 J				
	BENZO(B)FLUORANTHENE BENZO(G,H,I)PERYLENE	ug/kg ug/kg	4400 J 2400 J	6600 J 2900 J	12 U 15 U	300 J 220 J				
	BENZO(K)FLUORANTHENE	ug/kg ug/kg	1700	2200	17 U	19 U				
	BENZYL BUTYL PHTHALATE	ug/kg	20 U	19 U	18 U	19 U				
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	7000	11000	14 U	15 U				
	BIS(2-CHLOROETHOXY) METHANE	ug/kg	24 U	23 U	21 U	23 U				
	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/kg	20 U	19 U	18 U	19 U				
	BIS(2-CHLOROISOPROPYL) ETHER BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	17 U 15 U	17 U 14 U	15 U	17 U 14 U				
	CAPROLACTAM	ug/kg ug/kg	19 U	14 U 19 U	13 U 17 U	14 U				
	CARBAZOLE	ug/kg ug/kg	550	700	8 U	8.9 U				
218-01-9	CHRYSENE	ug/kg	8000	12000	17 U	290 J				
	DIBENZ(A,H)ANTHRACENE	ug/kg	720	880	11 U	12 U				
	DIBENZOFURAN	ug/kg	2200	2800	14 U	16 U				
	DIMETHYL DITTIAL ATE	ug/kg	6.4 U	6.2 U	5.7 U	6.3 U				
	DIMETHYL PHTHALATE DI-N-BUTYL PHTHALATE	ug/kg ug/kg	221 J 32 U	226 J 31 U	9.9 U 29 U	215 J 32 U				
	DI-N-OCTYLPHTHALATE	ug/kg ug/kg	4.7 U	4.6 U	4.2 U	4.6 U				
	FLUORANTHENE	ug/kg	14000	23000	7.3 U	570				
86-73-7	FLUORENE	ug/kg	22000	27000	14 U	15 U				
	HEXACHLOROBENZENE	ug/kg	17 U	16 U	15 U	17 U				
	HEXACHLOROBUTADIENE	ug/kg	15 U	15 U	13 U	15 U				
	HEXACHLOROCYCLOPENTADIENE HEXACHLOROETHANE	ug/kg	10 U 18 U	9.7 U 18 U	8.9 U 16 U	9.8 U 18 U				
	INDENO(1,2,3-C,D)PYRENE	ug/kg ug/kg	2100	2700	16 U 12 U	18 U 170 J				
	ISOPHORONE	ug/kg ug/kg	14 U	13 U	12 U	170 J				
	NAPHTHALENE	ug/kg	220000	240000	13 U	14 U				
98-95-3	NITROBENZENE	ug/kg	16 U	15 U	14 U	15 U				
	N-NITROSODI-N-PROPYLAMINE	ug/kg	21 U	20 U	18 U	20 U				
	N-NITROSODIPHENYLAMINE	ug/kg	9.9 U	9.6 U	8.8 U	9.7 U				
	PENTACHLOROPHENOL PHENANTHRENE	ug/kg	28 U 60000	27 U 79000	25 U 9.9 U	28 U 480				
02-01-0		ug/kg			9.9 U 8.4 U					
	PHENOL	ug/kg	9.5 U	9.2 U	X 4 11	9.4 U				

				Dup of MW-7(36-38)		
Con Ed - Luc	llow Site	Location ID:	MW- 7	MW- 7	MW- 7	MW-8
Validated So	l Analytical Data	Sample ID:	MW-7(36-38)	MW-17(36-38)	MW-7(48-50)	MW-8(9.5-10)
SDGs: D4710), D1751, D5300	Lab Sample Id:	D4710-03	D4710-09	D4710-04	D4751-03
		Depth:	36 - 38 ft	36 - 38 ft	48 - 50 ft	9.5 - 10 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4710	D4710	D4710	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	10/26/2012 9:10	10/26/2012 9:15	10/26/2012 9:20	11/6/2012 14:15
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	7150	6870	2050	4920 J
7440-36-0	ANTIMONY	mg/kg	0.28 U	0.28 U	0.25 U	0.28 UJ
7440-38-2	ARSENIC	mg/kg	2.45	2.47	1.55	9.24 J
7440-39-3	BARIUM	mg/kg	57.6	54	36.9	73 J
7440-41-7	BERYLLIUM	mg/kg	0.28	0.29	0.08 J	0.03 U
7440-43-9	CADMIUM	mg/kg	0.48	0.45	0.03 J	1.32
7440-70-2	CALCIUM	mg/kg	1350	1340	709	3220 J
7440-47-3	CHROMIUM, TOTAL	mg/kg	13.6 J	13.7 J	4.91 J	31.2 J
7440-48-4	COBALT	mg/kg	7.95	7.44	2.66	6.63
7440-50-8	COPPER	mg/kg	16 J	16.2 J	8.56 J	93.7 J
7439-89-6	IRON	mg/kg	22400	21300	5950	33500
7439-92-1	LEAD	mg/kg	14.4	14.6	3.07	170
7439-95-4	MAGNESIUM	mg/kg	3170	3100	1310	2310 J
7439-96-5	MANGANESE	mg/kg	180	135	44.7	240 J
7439-97-6	MERCURY	mg/kg	0.011 J	0.006 J	0.003 J	4.03
7440-02-0	NICKEL	mg/kg	23	22	8.83	19.4
7440-09-7	POTASSIUM	mg/kg	1980	1960	562	1530 J
7782-49-2	SELENIUM	mg/kg	0.21 U	0.21 U	0.52	0.57
7440-22-4	SILVER	mg/kg	0.5	0.47	0.1 J	0.36
7440-23-5	SODIUM	mg/kg	2090	2190	195	627 J
7440-28-0	THALLIUM	mg/kg	0.46 J	0.48 J	0.12 U	1.23
7440-62-2	VANADIUM	mg/kg	18	17.7	5.98	15.3 J
7440-66-6	ZINC	mg/kg	36.1	39.3	13.7	96.8
57-12-5	CYANIDE	mg/kg	0.29 J	0.362	0.171 J	0.097 J

Con Ed - Lud	low Site	Location ID:	MW-8	MW-9	MW-9	SB- 9
	Analytical Data	Sample ID:	MW-8(24.5-25)	MW-9(9-9.5)	MW-9(35-35.5)	SB-9(5-8)
SDGs: D4710	, D1751, D5300	Lab Sample Id:	D4751-04	D4751-06	D4751-05	D4710-07
		Depth:	24.5 - 25 ft	9 - 9.5 ft	35 - 35.5 ft	5 - 8 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4751	D4751	D4751	D4710
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	11/6/2012 14:20	11/6/2012 10:00	11/6/2012 10:05	10/26/2012 14:15
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	VOLATILES			0.07.77		0.00 **
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	1.1 U	0.95 U	0.99 U	0.99 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.59 U	0.5 U	0.52 U	0.52 UJ
76-13-1 79-00-5	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	1.7 U 1.2 U	1.4 U 0.97 U	1.5 U 1 U	1.5 U 1 UJ
75-34-3	1,1,2-TRICHLOROETHANE 1,1-DICHLOROETHANE	ug/kg ug/kg	1.2 U	0.97 U	1.1 U	1.1 U
75-34-3 75-35-4	1,1-DICHLOROETHANE 1,1-DICHLOROETHENE	ug/kg ug/kg	1.2 U 1.9 U	1.6 U	1.1 U 1.6 U	1.7 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg ug/kg	0.64 U	0.54 U	0.56 U	0.56 UJ
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg ug/kg	0.04 U	0.76 U	0.78 U	0.79 UJ
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg ug/kg	1.1 UJ	0.76 U 0.94 UJ	0.78 U 0.98 UJ	0.79 UJ 0.98 UJ
106-93-4	1,2-DIBROMO-5-CHLOROFROFANE 1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/kg ug/kg	0.82 U	0.69 U	0.72 U	0.72 UJ
95-50-1	1,2-DICHLOROBENZENE	ug/kg ug/kg	0.82 U 0.79 U	0.69 U 0.67 U	0.72 U	0.72 UJ
107-06-2	1,2-DICHLOROBENZENE 1,2-DICHLOROETHANE	ug/kg ug/kg	0.79 U	0.69 U	0.7 U	0.72 UJ
78-87-5	1,2-DICHLOROPROPANE	ug/kg ug/kg	0.32 U	0.28 U	0.72 U 0.29 U	0.72 UJ 0.29 UJ
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.47 U	0.4 U	0.41 U	0.42 UJ
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.53 U	0.44 U	0.46 U	0.46 UJ
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	64 U	54 U	55 U	56 UJ
591-78-6	2-HEXANONE	ug/kg	5 U	4.2 U	4.4 U	4.4 UJ
67-64-1	ACETONE	ug/kg	3.9 U	3.3 U	3.4 U	140
71-43-2	BENZENE	ug/kg	0.49 U	0.41 U	0.43 U	28 J
74-97-5	BROMOCHLOROMETHANE	ug/kg	1 U	0.85 U	0.89 U	0.89 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.79 U	0.67 U	0.7 U	0.7 UJ
75-25-2	BROMOFORM	ug/kg	0.95 U	0.8 U	0.83 U	0.83 UJ
74-83-9	BROMOMETHANE	ug/kg	3.1 U	2.6 U	2.7 U	2.8 U
75-15-0	CARBON DISULFIDE	ug/kg	1.4 U	1.1 U	1.2 U	1.2 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	1.3 U	1.1 U	1.1 U	1.1 UJ
108-90-7	CHLOROBENZENE	ug/kg	0.64 U	0.54 U	0.56 U	0.56 UJ
75-00-3	CHLOROETHANE	ug/kg	1.8 U	1.5 U	1.6 U	1.6 U
67-66-3	CHLOROFORM	ug/kg	0.95 U	0.8 U	0.83 U	0.83 U
74-87-3	CHLOROMETHANE	ug/kg	1.1 U	0.93 U	0.96 U	0.97 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	1.1 U	0.96 U	1 U	1 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.92 U	0.78 U	0.81 U	0.81 UJ
110-82-7	CYCLOHEXANE	ug/kg	1.3 U	1.1 U	1.1 U	1.1 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.69 U	0.58 U	0.61 U	0.61 UJ
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.83 U	0.7 U	0.73 U	0.73 U
100-41-4	ETHYLBENZENE	ug/kg	0.79 U	0.67 U	0.7 U	3.1 J
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.62 U	0.52 U	0.54 U	0.54 UJ
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.92 U	0.78 U	0.81 U	2.1 J
79-20-9	METHYL ACETATE	ug/kg	1.9 U	1.6 U	1.7 U	1.7 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/kg	4 U	3.4 U	3.5 U	3.5 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/kg	3.7 U	3.2 U	3.3 U	3.3 UJ
108-87-2	METHYLCYCLOHEXANE	ug/kg	1.4 U	1.1 U	1.2 U	1.2 UJ
75-09-2	METHYLENE CHLORIDE	ug/kg	11 UJ	5.8 UJ	5.6 UJ	50 UJ
95-47-6 100-42-5	O-XYLENE (1,2-DIMETHYLBENZENE) STYRENE	ug/kg	0.87 U 0.58 U	0.73 U 0.49 U	0.76 U 0.5 U	0.76 UJ 0.51 UJ
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.58 U 1.2 U	0.49 U 1 U	0.5 U 1.1 U	0.51 UJ 1.1 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	1.2 U 1.3 U	1.1 U	1.1 U 1.1 U	1.1 UJ
127-18-4	TOLUENE	ug/kg	0.82 U	0.69 U	0.72 U	1.1 UJ 11 J
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg ug/kg	0.82 U	0.69 U 0.75 U	0.72 U 0.77 U	0.78 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg ug/kg	0.88 U 1 U	0.75 U 0.85 U	0.77 U 0.89 U	0.78 U 0.89 UJ
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg ug/kg	1.1 U	0.83 U 0.93 U	0.89 U 0.96 U	0.89 UJ 0.97 UJ
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg ug/kg	1.1 U 1.7 U	0.93 U 1.4 U	1.5 U	0.97 UJ 1.5 U
75-09-4	VINYL CHLORIDE	ug/kg ug/kg	1.7 U 1.6 U	1.4 U	1.3 U 1.4 U	1.5 U 1.4 U
75 'U1- T	THE CHECKIPE	ug/Kg	1.0 0	1.5 0	1.7 0	1.7 U

Con Ed I "	vr. Cita	Logotion ID.	MW 0	MWA	MW	CD O
Con Ed - Ludlo	w Site Analytical Data	Location ID: Sample ID:	MW-8 MW-8(24.5-25)	MW-9 MW-9(9-9.5)	MW-9 MW-9(35-35.5)	SB- 9 SB-9(5-8)
SDGs: D4710,		Sample ID: Lab Sample Id:	MW-8(24.5-25) D4751-04	MW-9(9-9.5) D4751-06	MW-9(35-35.5) D4751-05	SB-9(5-8) D4710-07
3DG8. D4/10,	D1751, D5500	Depth:	24.5 - 25 ft	9 - 9.5 ft	35 - 35.5 ft	5 - 8 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4751	D4751	D4751	D4710
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	11/6/2012 14:20	11/6/2012 10:00	11/6/2012 10:05	10/26/2012 14:15
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
	SEMIVOLATILES	CIVIIS.				
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	17 U	14 U	15 U	15 U
	2,3,4,6-TETRACHLOROPHENOL	ug/kg	17 U	14 U	15 U	15 U
	2,4,5-TRICHLOROPHENOL	ug/kg	30 U	25 U	26 U	26 U
	2,4,6-TRICHLOROPHENOL	ug/kg	13 U	11 U	11 U	11 U
	2,4-DICHLOROPHENOL	ug/kg	16 U	14 U	14 U	14 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	24 U	21 U	21 U	21 U
51-28-5	2,4-DINITROPHENOL	ug/kg	43 UJ	37 UJ	38 UJ	38 U
	2,4-DINITROTOLUENE	ug/kg	13 U	11 U	11 U	11 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	17 U	15 U	15 U	15 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	9.7 U	8.2 U	8.5 U	8.5 U
95-57-8	2-CHLOROPHENOL	ug/kg	22 U	19 U	20 U	20 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	11 U	9.1 U	9.4 U	9.4 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	23 U	20 U	20 U	20 U
	2-NITROANILINE	ug/kg	19 U	16 U	17 U	17 U
88-75-5	2-NITROPHENOL	ug/kg	21 U	17 U	18 U	18 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	27 U	23 U	24 U	24 U
МЕРНЗМЕРН	3+4-Methylphenols	ug/kg	22 U	19 U	19 U	19 U
99-09-2	3-NITROANILINE	ug/kg	27 U	23 U	24 U	24 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	24 U	21 UJ	21 U	21 U
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	8.3 U	7.1 U	7.3 U	7.3 U
	4-CHLORO-3-METHYLPHENOL	ug/kg	19 U	16 U	17 U	17 U
	4-CHLOROANILINE	ug/kg	30 UJ	26 UJ	26 UJ	26 U
	4-CHLOROPHENYL PHENYL ETHER	ug/kg	23 U	20 U	20 U	20 U
	4-NITROANILINE	ug/kg	55 U	47 U	49 U	49 U
	4-NITROPHENOL	ug/kg	79 U	67 U	69 U	69 U
83-32-9	ACENAPHTHENE	ug/kg	12 U	10 U	11 U	11 U
208-96-8	ACENAPHTHYLENE	ug/kg	11 U	9.1 U	9.4 U	9.4 U
	ACETOPHENONE	ug/kg	13 U	11 U	11 U	11 U
	ANTHRACENE	ug/kg	8.7 U	7.4 U	7.6 U	7.6 U
	ATRAZINE	ug/kg	22 U	19 U	20 U	20 U
	BENZALDEHYDE	ug/kg	22 U	19 U	19 U	20 U
	BENZO(A)ANTHRACENE	ug/kg	20 U	17 U	18 U	290 Ј
	BENZO(A)PYRENE	ug/kg	9.2 U	7.8 U	8.1 U	280 J
	BENZO(B)FLUORANTHENE	ug/kg	14 U	12 U	12 U	380 J
	BENZO(G,H,I)PERYLENE	ug/kg	17 U	15 U	15 U	170 J
	BENZO(K)FLUORANTHENE	ug/kg	20 U	17 U	18 U	18 U
	BENZYL BUTYL PHTHALATE	ug/kg	20 U	17 U	18 U	18 U
	BIPHENYL (DIPHENYL)	ug/kg	16 U	14 U	14 U	14 U
	BIS(2-CHLOROETHOXY) METHANE	ug/kg	25 U	21 U	22 U	22 U
	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/kg	20 U	17 U	18 U	18 U
	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	18 U	15 U	15 U	15 U
	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	15 U	13 U	13 U	13 U
105-60-2	CAPROLACTAM	ug/kg	20 U	17 U	17 U	17 U
	CARBAZOLE	ug/kg	9.3 U	7.9 U	8.2 U	8.2 U
	CHRYSENE DIRENZ(A L) ANTURACENIE	ug/kg	19 U	16 U	17 U	320 J
	DIBENZ(A,H)ANTHRACENE DIBENZOFURAN	ug/kg	12 U 17 U	10 U 14 U	11 U 15 U	11 U 15 U
		ug/kg				
	DIETHYL PHTHALATE	ug/kg	6.6 U	5.6 U	5.8 U	5.8 U
	DIMETHYL PHTHALATE	ug/kg	221 J	292 J	200 J	161 J
	DI-N-BUTYL PHTHALATE DI-N-OCTYLPHTHALATE	ug/kg	33 U	28 U	29 U	29 U
		ug/kg	4.9 U	4.1 U 7.3 U	4.3 U 7.5 U	4.3 U 560
	FLUORANTHENE ELHOPENE	ug/kg	8.6 U	7.3 U 14 U	7.5 U 14 U	360 14 U
	FLUORENE HEXACHLOROBENZENE	ug/kg	16 U 17 U	14 U 15 U	14 U 15 U	14 U 15 U
	HEXACHLOROBENZENE HEXACHLOROBUTADIENE	ug/kg	17 U 15 U	15 U 13 U	15 U 14 U	15 U 14 U
	HEXACHLOROCYCLOPENTADIENE	ug/kg	10 U	8.8 U	9.1 U	9.1 U
	HEXACHLOROETHANE	ug/kg ug/kg	10 U	8.8 U 16 U	9.1 U 17 U	9.1 U 17 U
	INDENO(1,2,3-C,D)PYRENE	ug/kg ug/kg	19 U 14 U	16 U 12 U	17 U 12 U	17 U 190 J
	INDENO(1,2,3-C,D)PYRENE ISOPHORONE		14 U 14 U	12 U 12 U	12 U 12 U	190 J 12 U
	NAPHTHALENE	ug/kg	14 U 15 U	12 U 12 U	12 U 13 U	12 U 13 U
	NITROBENZENE	ug/kg ug/kg	16 U	12 U 14 U	13 U 14 U	13 U 14 U
	N-NITROSODI-N-PROPYLAMINE	ug/kg ug/kg	21 U	18 U	14 U 19 U	14 U
	N-NITROSODI-N-PROPTLAMINE N-NITROSODIPHENYLAMINE	ug/kg ug/kg	10 U	8.7 U	9 U	9 U
	PENTACHLOROPHENOL	ug/kg ug/kg	29 U	25 U	26 U	26 U
	PHENANTHRENE	ug/kg ug/kg	11 U	9.8 U	10 U	320 J
	PHENOL	ug/kg ug/kg	9.8 U	8.4 U	8.6 U	8.6 U
	PYRENE	ug/kg ug/kg	10 U	8.7 U	9 U	470
127-00-0	LINEAL	ug/Ng	10 0	0.7 U	7 0	470

Con Ed - Luc	llow Site	Location ID:	MW-8	MW-9	MW-9	SB- 9
Validated So	il Analytical Data	Sample ID:	MW-8(24.5-25)	MW-9(9-9.5)	MW-9(35-35.5)	SB-9(5-8)
SDGs: D471	0, D1751, D5300	Lab Sample Id:	D4751-04	D4751-06	D4751-05	D4710-07
		Depth:	24.5 - 25 ft	9 - 9.5 ft	35 - 35.5 ft	5 - 8 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4751	D4751	D4751	D4710
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	11/6/2012 14:20	11/6/2012 10:00	11/6/2012 10:05	10/26/2012 14:15
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	8550 J	13400 J	2710 J	7070
7440-36-0	ANTIMONY	mg/kg	0.3 UJ	0.26 UJ	0.26 UJ	0.26 U
7440-38-2	ARSENIC	mg/kg	2.16 J	0.15 U	0.75 J	7.78
7440-39-3	BARIUM	mg/kg	85.6 J	177 J	60.2 J	157
7440-41-7	BERYLLIUM	mg/kg	0.33	0.13 J	0.07 J	0.11 J
7440-43-9	CADMIUM	mg/kg	0.66	2.17	0.19	1.24
7440-70-2	CALCIUM	mg/kg	1870 J	2660 J	13000 J	5480
7440-47-3	CHROMIUM, TOTAL	mg/kg	15.8 J	28.1 J	6.01 J	16.7 J
7440-48-4	COBALT	mg/kg	10.7	27.8	3.85	7.34
7440-50-8	COPPER	mg/kg	18.8 J	7.84 J	8.4 J	44.1 J
7439-89-6	IRON	mg/kg	24400	41900	6990	29900
7439-92-1	LEAD	mg/kg	16.1	17.4	4.01	462
7439-95-4	MAGNESIUM	mg/kg	4040 J	7840 J	6590 J	2280
7439-96-5	MANGANESE	mg/kg	427 J	251 J	192 J	284
7439-97-6	MERCURY	mg/kg	0.002 U	0.425	0.002 U	2.33
7440-02-0	NICKEL	mg/kg	26.5	48.6	10.6	21.2
7440-09-7	POTASSIUM	mg/kg	2600 J	10700 J	860 J	655
7782-49-2	SELENIUM	mg/kg	0.26 J	0.5	0.19 U	0.49
7440-22-4	SILVER	mg/kg	0.57	0.23 J	0.07 U	0.79
7440-23-5	SODIUM	mg/kg	662 J	402 J	1080 J	1150
7440-28-0	THALLIUM	mg/kg	0.52 J	0.12 U	0.12 U	1.13
7440-62-2	VANADIUM	mg/kg	21.2 J	38.8 J	9.23 J	16.7
7440-66-6	ZINC	mg/kg	47.2	97.6	14.5	268
57-12-5	CYANIDE	mg/kg	0.042 U	0.087 J	0.075 J	0.211 J

G EL 1 11	ar.	Tr .: 175	ar o	GD 10	GD 10	ap 11
Con Ed - Ludl		Location ID:	SB- 9	SB-10	SB-10	SB-11
	Analytical Data	Sample ID:	SB-9(28-30)	SB-10(30-32)	SB-10(45-47)	SB-11(5-5.5)
SDGs: D4/10,	, D1751, D5300	Lab Sample Id:	D4710-08	D4710-01	D4710-02	D4751-01
		Depth:	28 - 30 ft	30 - 32 ft	45 - 47 ft	5 - 5.5 ft
		Source: SDG:	CTECH D4710	CTECH D4710	CTECH D4710	CTECH D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	10/26/2012 14:55	10/25/2012 12:10	10/25/2012 12:40	11/6/2012 12:05
					1/23/2012 12:40	
CAS NO.	COMPOUND	Validated: UNITS:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	VOLATILES	UNIIS:				
71-55-6	1,1,1-TRICHLOROETHANE	na/ka	1.1 U	5.3 U	1.1 U	1.2 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.6 U	3.3 U R	0.55 U	0.64 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	1.7 U	8 U	1.6 U	1.8 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg ug/kg	1.7 U	5.4 U	1.0 U	1.8 U
75-34-3	1,1-DICHLOROETHANE		1.2 U	5.7 U	1.1 U	1.2 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	1.2 U 1.9 U	8.9 U	1.1 U 1.8 U	2 U
73-33-4 87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg				
120-82-1		ug/kg	0.65 U	R	0.6 U	0.69 U
	1,2,4-TRICHLOROBENZENE	ug/kg	0.91 U	R	0.84 U	0.97 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1.1 U	R	1 U	1.2 UJ
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/kg	0.83 U	3.9 U	0.77 U	0.89 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.81 U	R	0.75 U	0.86 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.83 U	3.9 U	0.77 U	0.89 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.34 U	1.6 U	0.31 U	0.36 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.48 U	R	0.45 U	0.51 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.53 U	R	0.49 U	0.57 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	65 U	300 U	60 U	69 U
591-78-6	2-HEXANONE	ug/kg	5.1 U	24 U	4.7 U	5.4 U
67-64-1	ACETONE	ug/kg	17 J	18 U	3.6 U	4.2 U
71-43-2	BENZENE	ug/kg	7.9	210000	0.46 U	0.53 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	1 U	4.8 U	0.95 U	1.1 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.81 U	3.7 U	0.75 U	0.86 U
75-25-2	BROMOFORM	ug/kg	0.96 U	4.5 U	0.89 U	1 U
74-83-9	BROMOMETHANE	ug/kg	3.2 U	15 U	3 U	3.4 U
75-15-0	CARBON DISULFIDE	ug/kg	1.4 U	41	1.3 U	2.8 J
56-23-5	CARBON TETRACHLORIDE	ug/kg	1.3 U	6 U	1.2 U	1.4 U
108-90-7	CHLOROBENZENE	ug/kg	0.65 U	3 U	0.6 U	0.69 U
75-00-3	CHLOROETHANE	ug/kg	1.8 U	8.5 U	1.7 U	1.9 U
67-66-3	CHLOROFORM	ug/kg	0.96 U	4.5 U	0.89 U	1 U
74-87-3	CHLOROMETHANE	ug/kg	1.1 U	5.2 U	1 U	1.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	1.2 U	5.4 U	1.1 U	1.2 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.94 U	4.3 U	0.87 U	1 U
110-82-7	CYCLOHEXANE	ug/kg	1.3 U	57	1.2 U	1.4 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.7 U	3.3 U	0.65 U	0.75 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.84 U	3.9 U	0.78 U	0.9 U
100-41-4	ETHYLBENZENE	ug/kg	2 J	540000	0.75 U	0.86 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.62 U	29000 J	0.58 U	0.67 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	1.5 J	470000	0.87 U	1 U
79-20-9 78-93-3	METHYL ACETATE	ug/kg	2 U	9.1 U	1.8 U	2.1 U
	METHYL ETHYL KETONE (2-BUTANONE)	ug/kg	4 U	19 U	3.7 U	4.3 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/kg	3.8 U	18 U	3.5 U	4.1 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	1.4 U	180	1.3 U	1.5 U
75-09-2 95-47-6	METHYLENE CHLORIDE	ug/kg	25 UJ	160 J	7.1 UJ	8.7 UJ
95-47-6 100-42-5	O-XYLENE (1,2-DIMETHYLBENZENE) STYRENE	ug/kg	0.88 U 0.58 U	220000 7100 J	0.82 U 0.54 U	0.94 U 0.62 U
1634-04-4		ug/kg				
	TERT-BUTYL METHYL ETHER	ug/kg	1.2 U	5.8 U	1.2 U	1.3 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	1.3 U	6.1 U	1.2 U	1.4 U
108-88-3	TOLUENE TRANS 1.2 DICHI OPOETHENE	ug/kg	5.9 J	510000	0.77 U	0.89 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.9 U	4.2 U	0.83 U	0.96 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	1 U	4.8 U	0.95 U	1.1 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	1.1 U	5.2 U	1 U	1.2 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg	1.7 U	8 U	1.6 U	1.8 U
75-01-4	VINYL CHLORIDE	ug/kg	1.6 U	7.4 U	1.5 U	1.7 U

C F1 I II	C'A	Ir .: ID	CD 0	CD 10	CD 10	CD 11
Con Ed - Ludlo		Location ID:	SB- 9	SB-10	SB-10	SB-11
	Analytical Data	Sample ID: Lab Sample Id:	SB-9(28-30) D4710-08	SB-10(30-32)	SB-10(45-47)	SB-11(5-5.5)
SDGs: D4710,	D1/51, D5500			D4710-01	D4710-02	D4751-01
		Depth:	28 - 30 ft	30 - 32 ft	45 - 47 ft	5 - 5.5 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4710	D4710	D4710	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	10/26/2012 14:55	10/25/2012 12:10	10/25/2012 12:40	11/6/2012 12:05
~.~.	gor movne	Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES			100 **		40.77
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	17 U	480 U	16 U	18 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	17 U	480 U	16 U	18 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	30 U	850 U	28 U	32 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	13 U	370 U	12 U	14 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	16 U	460 U	15 U	18 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	25 U	690 U	23 U	26 U
51-28-5	2,4-DINITROPHENOL	ug/kg	44 U	1200 U	41 U	47 UJ
121-14-2	2,4-DINITROTOLUENE	ug/kg	13 U	370 U	12 U	14 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	18 U	500 U	16 U	19 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	9.9 U	280 U	9.1 U	11 U
95-57-8	2-CHLOROPHENOL	ug/kg	23 U	640 U	21 U	24 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	11 U	2800000	10 U	12 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	23 U	660 U	22 U	25 U
88-74-4	2-NITROANILINE	ug/kg	19 U	540 U	18 U	21 U
88-75-5	2-NITROPHENOL	ug/kg	21 U	590 U	19 U	22 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	28 U	780 U	26 U	30 U
	3+4-Methylphenols	ug/kg	22 U	630 U	21 U	24 U
99-09-2	3-NITROANILINE	ug/kg	28 U	780 U	26 U	30 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	25 U	700 U	23 U	27 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	8.4 U	240 U	7.8 U	9 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	19 U	540 U	18 U	21 U
106-47-8	4-CHLOROANILINE	ug/kg	30 U	860 U	28 U	33 UJ
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	23 U	660 U	22 U	25 U
100-01-6	4-NITROANILINE	ug/kg	56 U	1600 U	52 U	60 U
100-02-7	4-NITROPHENOL	ug/kg	80 U	2300 U	74 U	86 U
83-32-9	ACENAPHTHENE	ug/kg	12 U	320000 J	11 U	13 U
208-96-8	ACENAPHTHYLENE	ug/kg	11 U	370000	200 J	12 U
98-86-2	ACETOPHENONE	ug/kg	13 U	370 U	12 U	14 U
120-12-7	ANTHRACENE	ug/kg	8.8 U	370000	8.2 U	9.4 U
1912-24-9	ATRAZINE	ug/kg	23 U	640 U	21 U	24 U
100-52-7	BENZALDEHYDE	ug/kg	23 U	640 U	21 U	24 U
56-55-3	BENZO(A)ANTHRACENE	ug/kg	21 U	220000	19 U	22 U
50-32-8	BENZO(A)PYRENE	ug/kg	9.3 U	150000 J	8.7 U	10 U
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	14 U	97000 J	13 U	15 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	18 U	42000 J	16 U	19 U
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	20 U	32000	19 U	22 U
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	21 U	580 U	19 U	22 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	16 U	200000	15 U	17 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	25 U	700 U	23 U	27 U
111-44-4	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/kg	21 U	580 U	19 U	22 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	18 U	500 U	17 U	19 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	15 U	430 U	14 U	16 U
105-60-2	CAPROLACTAM	ug/kg	20 U	570 U	19 U	22 U
86-74-8	CARBAZOLE	ug/kg	9.5 U	13000	8.8 U	10 U
218-01-9	CHRYSENE	ug/kg	20 U	220000	18 U	21 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	12 U	12000	12 U	13 U
132-64-9	DIBENZOFURAN	ug/kg	17 U	43000	16 U	18 U
84-66-2	DIETHYL PHTHALATE	ug/kg	6.7 U	190 U	6.3 U	7.2 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	12 U	330 U	11 U	296 J
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	34 U	960 U	32 U	36 U
117-84-0	DI-N-OCTYLPHTHALATE	ug/kg	4.9 U	140 U	4.6 U	5.3 U
206-44-0	FLUORANTHENE	ug/kg	8.7 U	400000	8.1 U	9.3 U
86-73-7	FLUORENE	ug/kg	16 U	550000	15 U	17 U
118-74-1	HEXACHLOROBENZENE	ug/kg	18 U	500 U	16 U	19 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	16 U	440 U	15 U	17 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	11 U	300 U	9.7 U	11 U
67-72-1	HEXACHLOROETHANE	ug/kg	19 U	540 U	18 U	21 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	14 U	37000	13 U	15 U
78-59-1	ISOPHORONE	ug/kg	14 U	400 U	13 U	15 U
91-20-3	NAPHTHALENE	ug/kg	15 U	5700000	210 J	16 U
98-95-3	NITROBENZENE	ug/kg	16 U	460 U	15 U	17 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	22 U	610 U	20 U	23 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	10 U	290 U	9.6 U	11 U
87-86-5	PENTACHLOROPHENOL	ug/kg	30 U	830 U	27 U	32 U
85-01-8	PHENANTHRENE	ug/kg	12 U	1500000	11 U	12 U
108-95-2 129-00-0	PHENOL	ug/kg	10 U	280 U	9.3 U	11 U
	PYRENE	ug/kg	10 U	640000	9.6 U	11 U

Con Ed - Luc	llow Site	Location ID:	SB- 9	SB-10	SB-10	SB-11
Validated So	il Analytical Data	Sample ID:	SB-9(28-30)	SB-10(30-32)	SB-10(45-47)	SB-11(5-5.5)
SDGs: D471	0, D1751, D5300	Lab Sample Id:	D4710-08	D4710-01	D4710-02	D4751-01
		Depth:	28 - 30 ft	30 - 32 ft	45 - 47 ft	5 - 5.5 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4710	D4710	D4710	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	10/26/2012 14:55	10/25/2012 12:10	10/25/2012 12:40	11/6/2012 12:05
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	9490	3130	4490	4170 J
7440-36-0	ANTIMONY	mg/kg	0.31 U	0.28 U	0.29 U	0.34 UJ
7440-38-2	ARSENIC	mg/kg	3	7.21	1.23	18.3 J
7440-39-3	BARIUM	mg/kg	96	30.9	55.6	156 J
7440-41-7	BERYLLIUM	mg/kg	0.39	0.04 J	0.2	0.04 U
7440-43-9	CADMIUM	mg/kg	0.7	1.75	0.29	0.16 J
7440-70-2	CALCIUM	mg/kg	2540	9540	13000	2260 J
7440-47-3	CHROMIUM, TOTAL	mg/kg	18.4 J	5.38 J	8.9 J	9.93 J
7440-48-4	COBALT	mg/kg	11.9	3.4	5.38	6.31
7440-50-8	COPPER	mg/kg	20.5 J	26.5 J	11.8 J	69.6 J
7439-89-6	IRON	mg/kg	28600	13200	12600	26800
7439-92-1	LEAD	mg/kg	18.1	215	8.81	30
7439-95-4	MAGNESIUM	mg/kg	5050	5480	5450	1690 J
7439-96-5	MANGANESE	mg/kg	414	277	290	803 J
7439-97-6	MERCURY	mg/kg	0.002 U	0.381	0.002 U	1.17
7440-02-0	NICKEL	mg/kg	31.1	12.2	14.3	17.5
7440-09-7	POTASSIUM	mg/kg	2730	352	1470	1500 J
7782-49-2	SELENIUM	mg/kg	0.33 J	1.95	0.21 U	1
7440-22-4	SILVER	mg/kg	0.67	0.15 J	0.09 J	0.51
7440-23-5	SODIUM	mg/kg	148	4300	150	867 J
7440-28-0	THALLIUM	mg/kg	0.74 J	1.15	0.18 J	0.46 J
7440-62-2	VANADIUM	mg/kg	22.5	7.59	13.1	14.8 J
7440-66-6	ZINC	mg/kg	50.8	506	22.9	88.4
57-12-5	CYANIDE	mg/kg	0.244 J	4.4	0.172 J	0.133 J

Con Ed - Ludle	ow Site	Location ID:	SB-11	SB-12	SB-12	SB-13
	Analytical Data	Sample ID:	SB-11(29.5-30)	SB-12(12-14)	SB-12(36-38)	SB-13(47-49)
SDGs: D4710.	, D1751, D5300	Lab Sample Id:	D4751-02	D4751-07	D4751-08	D4751-09
		Depth:	29.5 - 30 ft	12 - 14 ft	36 - 38 ft	47 - 49 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4751	D4751	D4751	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	11/6/2012 12:10	11/5/2012 10:10	11/5/2012 11:00	11/5/2012 15:05
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	1.1 U	44 U	0.93 U	1 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.55 U	34 U	0.48 U	0.54 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	1.6 U	50 U	1.4 U	1.6 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.1 U	42 U 40 U	0.95 U	1.1 U 1.1 U
75-34-3 75-35-4	1,1-DICHLOROETHANE	ug/kg	1.1 U 1.8 U	52 U	0.99 U 1.5 U	1.1 U 1.7 U
73-33-4 87-61-6	1,1-DICHLOROETHENE	ug/kg				
120-82-1	1,2,3-TRICHLOROBENZENE 1,2,4-TRICHLOROBENZENE	ug/kg	0.6 U 0.84 U	72 U 69 U	0.53 U 0.74 U	0.58 U 0.82 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	0.84 U 1 U	51 U	0.74 U 0.92 UJ	0.82 U 1 UJ
106-93-4	1,2-DIBROMO-3-CHLOROPROPANE 1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/kg ug/kg	0.77 U	46 U	0.92 UJ 0.67 U	0.75 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg ug/kg	0.77 U	50 U	0.65 U	0.73 U
107-06-2	1,2-DICHLOROETHANE	ug/kg ug/kg	0.77 U	53 U	0.67 U	0.75 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg ug/kg	0.77 U	51 U	0.07 U	0.73 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.45 U	48 U	0.39 U	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.49 U	36 U	0.43 U	0.48 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	60 U	5500 U	53 U	58 U
591-78-6	2-HEXANONE	ug/kg	4.7 U	220 UJ	4.1 U	4.6 U
67-64-1	ACETONE	ug/kg	3.6 UJ	340 UJ	3.2 U	3.5 U
71-43-2	BENZENE	ug/kg	0.46 U	36 U	0.4 U	85
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.95 U	250 U	0.83 U	0.92 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.75 U	40 U	0.65 U	0.73 U
75-25-2	BROMOFORM	ug/kg	0.89 U	52 U	0.78 U	0.87 U
74-83-9	BROMOMETHANE	ug/kg	3 U	69 U	2.6 U	2.9 U
75-15-0	CARBON DISULFIDE	ug/kg	1.3 U	60 U	1.1 U	2.6 J
56-23-5	CARBON TETRACHLORIDE	ug/kg	1.2 U	69 U	1 U	1.2 U
108-90-7	CHLOROBENZENE	ug/kg	0.6 U	54 U	0.53 U	0.58 U
75-00-3	CHLOROETHANE	ug/kg	1.7 U	73 UJ	1.5 U	1.6 U
67-66-3	CHLOROFORM	ug/kg	0.89 U	38 U	0.78 U	0.87 U
74-87-3	CHLOROMETHANE	ug/kg	1 U	60 U	0.91 U	1 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	1.1 U	39 U	0.94 U	1 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.87 U	34 U	0.76 U	0.84 U
110-82-7	CYCLOHEXANE	ug/kg	1.2 U	61 U	1.1 U	1.7 J
124-48-1 75-71-8	DIBROMOCHLOROMETHANE	ug/kg	0.65 U	58 U	0.57 U	0.63 U
100-41-4	DICHLORODIFLUOROMETHANE ETHYLBENZENE	ug/kg	0.78 U 0.75 U	61 U 9100	0.68 U 0.65 U	0.76 U 460 J
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.73 U 0.58 U	1000	0.65 U 0.51 U	460 J 45
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg ug/kg	0.87 U	11000	0.76 U	230
79-20-9	METHYL ACETATE	ug/kg ug/kg	1.8 U	92 UJ	1.6 U	1.8 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/kg ug/kg	3.7 U	150 U	3.3 U	3.6 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/kg ug/kg	3.5 U	230 U	3.1 U	3.4 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	1.3 U	69000	1.1 U	7.1
75-09-2	METHYLENE CHLORIDE	ug/kg	6 UJ	46 U	5.3 UJ	10 UJ
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.82 U	5700	0.72 U	160
100-42-5	STYRENE	ug/kg	0.54 U	290 J	0.47 U	3.8 J
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	1.2 U	39 U	1 U	1.1 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	1.2 U	30 U	1.1 U	1.2 U
108-88-3	TOLUENE	ug/kg	0.77 U	1500	0.67 U	53
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.83 U	46 U	0.73 U	0.81 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.95 U	32 U	0.83 U	0.92 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	1 U	31 U	0.91 U	1 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg	1.6 U	39 U	1.4 U	1.5 U
75-01-4	VINYL CHLORIDE	ug/kg	1.5 U	38 U	1.3 U	1.4 U

C E1 111-	C:4-	Location ID:	CD 11	CD 12	CD 12	CD 12
	Con Ed - Ludlow Site Validated Soil Analytical Data		SB-11	SB-12	SB-12	SB-13
			SB-11(29.5-30)	SB-12(12-14)	SB-12(36-38)	SB-13(47-49)
SDGs: D4/10,	D1/51, D5300	Lab Sample Id:	D4751-02	D4751-07	D4751-08	D4751-09
		Depth:	29.5 - 30 ft	12 - 14 ft	36 - 38 ft	47 - 49 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4751	D4751	D4751	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	11/6/2012 12:10	11/5/2012 10:10	11/5/2012 11:00	11/5/2012 15:05
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
	COMPOUND	UNITS:				
	SEMIVOLATILES	<u>,</u>				
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	16 U	73 U	14 U	15 U
	2,3,4,6-TETRACHLOROPHENOL	ug/kg	16 U	73 U	14 U	15 U
	2,4,5-TRICHLOROPHENOL	ug/kg	28 U	130 U	25 U	28 U
	2,4,6-TRICHLOROPHENOL	ug/kg	12 U	57 U	11 U	12 U
	2,4-DICHLOROPHENOL	ug/kg	15 U	70 U	13 U	15 U
	2,4-DIMETHYLPHENOL	ug/kg	23 U	100 U	20 U	22 U
	2,4-DINITROPHENOL	ug/kg	41 UJ	190 UJ	36 UJ	40 UJ
	2,4-DINITROTOLUENE	ug/kg	12 U	56 U	11 U	12 U
	2,6-DINITROTOLUENE	ug/kg	16 U	75 U	14 U	16 U
	2-CHLORONAPHTHALENE	ug/kg	9.1 U	42 U	8 U	8.9 U
	2-CHLOROPHENOL	ug/kg	21 U	98 U	19 U	21 U
	2-METHYLNAPHTHALENE	ug/kg	10 U	160000	8.8 U	1300
	2-METHYLPHENOL (O-CRESOL)	ug/kg	22 U	100 U	19 U	21 U
	2-NITROANILINE	ug/kg	18 U	82 U	16 U	17 U
	2-NITROPHENOL	ug/kg	19 U	89 U	17 U	19 U
	3,3'-DICHLOROBENZIDINE	ug/kg	26 U	120 U	23 U	25 U
МЕРН3МЕРН	3+4-Methylphenols	ug/kg	21 U	96 U	18 U	20 U
	3-NITROANILINE	ug/kg	26 U	120 U	23 U	25 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	23 U	110 UJ	20 UJ	22 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	7.8 U	36 U	6.8 U	7.6 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	18 U	82 U	16 U	17 U
106-47-8	4-CHLOROANILINE	ug/kg	28 UJ	130 UJ	25 UJ	28 UJ
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	22 U	100 U	19 U	21 U
100-01-6	4-NITROANILINE	ug/kg	52 U	240 U	46 U	51 U
100-02-7	4-NITROPHENOL	ug/kg	74 U	340 U	65 U	73 U
83-32-9	ACENAPHTHENE	ug/kg	11 U	9100	9.9 U	320 J
208-96-8	ACENAPHTHYLENE	ug/kg	10 U	50000	8.8 U	240 J
98-86-2	ACETOPHENONE	ug/kg	12 U	57 U	11 U	12 U
120-12-7	ANTHRACENE	ug/kg	8.2 U	22000	7.2 U	310 J
1912-24-9	ATRAZINE	ug/kg	21 U	98 U	19 U	21 U
100-52-7	BENZALDEHYDE	ug/kg	21 U	96 U	18 U	20 U
56-55-3	BENZO(A)ANTHRACENE	ug/kg	19 U	11000	17 U	260 J
50-32-8	BENZO(A)PYRENE	ug/kg	8.6 U	7200	7.6 U	200 Ј
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	13 U	5500	11 U	160 J
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	16 U	2100	14 U	16 U
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	19 U	1700 J	17 U	18 U
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	19 U	89 U	17 U	19 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	15 U	10000	13 U	15 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	23 U	110 U	20 U	23 U
111-44-4	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/kg	19 U	89 U	17 U	19 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	17 U	77 U	15 U	16 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	14 U	65 U	12 U	14 U
105-60-2	CAPROLACTAM	ug/kg	19 U	86 U	16 U	18 U
	CARBAZOLE	ug/kg	8.8 U	40 U	7.7 U	8.6 U
	CHRYSENE	ug/kg	18 U	10000	16 U	260 J
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	12 U	53 U	10 U	11 U
132-64-9	DIBENZOFURAN	ug/kg	16 U	2500	14 U	15 U
84-66-2	DIETHYL PHTHALATE	ug/kg	6.2 U	29 U	5.5 U	6.1 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	284 J	50 U	455	403
	DI-N-BUTYL PHTHALATE	ug/kg	31 U	150 U	28 U	31 U
	DI-N-OCTYLPHTHALATE	ug/kg	4.6 U	21 U	4 U	4.5 U
	FLUORANTHENE	ug/kg	8 U	22000	7 U	490
	FLUORENE	ug/kg	15 U	33000	13 U	320 J
	HEXACHLOROBENZENE	ug/kg	16 U	75 U	14 U	16 U
	HEXACHLOROBUTADIENE	ug/kg	15 U	67 U	13 U	14 U
	HEXACHLOROCYCLOPENTADIENE	ug/kg	9.7 U	45 U	8.5 U	9.5 U
	HEXACHLOROETHANE	ug/kg	18 U	83 U	16 U	18 U
	INDENO(1,2,3-C,D)PYRENE	ug/kg	13 U	2000	12 U	13 U
	ISOPHORONE	ug/kg	13 U	61 U	12 U	13 U
	NAPHTHALENE	ug/kg	14 U	310000	12 U	2800
	NITROBENZENE	ug/kg	15 U	70 U	13 U	15 U
	N-NITROSODI-N-PROPYLAMINE	ug/kg	20 U	93 U	18 U	20 U
	N-NITROSODIPHENYLAMINE	ug/kg	9.6 U	44 U	8.4 U	9.4 U
	PENTACHLOROPHENOL	ug/kg	27 U	130 U	24 U	27 U
	PHENANTHRENE	ug/kg	11 U	94000	9.5 U	1300
	PHENOL	ug/kg	9.2 U	43 U	8.1 U	9 U
	PYRENE	ug/kg	9.6 U	29000	8.4 U	620

Con Ed - Lud	llow Site	Location ID:	SB-11	SB-12	SB-12	SB-13
	il Analytical Data	Sample ID:	SB-11(29.5-30)	SB-12(12-14)	SB-12(36-38)	SB-13(47-49)
), D1751, D5300	Lab Sample Id:	D4751-02	D4751-07	D4751-08	D4751-09
SDGS: D4/10	0, D1751, D5500		29.5 - 30 ft	12 - 14 ft	36 - 38 ft	47 - 49 ft
		Depth: Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4751	D4751	D4751	D4751
		SDG: Matrix:		SOIL		SOIL
			SOIL		SOIL	
		Sampled:	11/6/2012 12:10	11/5/2012 10:10	11/5/2012 11:00	11/5/2012 15:05
	T	Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	INORGANICS	4				
7429-90-5	ALUMINUM	mg/kg	5890 J	4120 J	2800 J	9270 J
7440-36-0	ANTIMONY	mg/kg	0.28 UJ	0.27 UJ	0.24 UJ	0.28 UJ
7440-38-2	ARSENIC	mg/kg	1.3 J	0.9 J	0.49 J	1.96 J
7440-39-3	BARIUM	mg/kg	63.3 J	31.9 J	19.1 J	124 J
7440-41-7	BERYLLIUM	mg/kg	0.3	0.14 J	0.04 J	0.47
7440-43-9	CADMIUM	mg/kg	0.42	0.35	0.18	0.71
7440-70-2	CALCIUM	mg/kg	12200 J	952 J	10100 J	19200 J
7440-47-3	CHROMIUM, TOTAL	mg/kg	11.7 J	9.3 J	7.04 J	19.9 J
7440-48-4	COBALT	mg/kg	6.51	6.88	3.39	7.91
7440-50-8	COPPER	mg/kg	12.1 J	11 J	9.25 J	18.3 J
7439-89-6	IRON	mg/kg	14600	10300	6990	18600
7439-92-1	LEAD	mg/kg	10.5	45.3	3.88	15.2
7439-95-4	MAGNESIUM	mg/kg	5740 J	1870 J	4920 J	3760 J
7439-96-5	MANGANESE	mg/kg	319 J	191 J	209 J	1240 J
7439-97-6	MERCURY	mg/kg	0.002 U	0.002 U	0.002 U	0.02
7440-02-0	NICKEL	mg/kg	16.6	11.2	11.5	21.4
7440-09-7	POTASSIUM	mg/kg	2190 J	957 J	581 J	2340 J
7782-49-2	SELENIUM	mg/kg	0.21 U	0.2 U	0.17 U	1.06
7440-22-4	SILVER	mg/kg	0.14 J	0.1 J	0.06 U	0.34
7440-23-5	SODIUM	mg/kg	2520 J	5360 J	1220 J	1280 J
7440-28-0	THALLIUM	mg/kg	0.34 J	0.13 U	0.13 J	0.81 J
7440-62-2	VANADIUM	mg/kg	15.3 J	12.7 J	9.49 J	18.3 J
7440-66-6	ZINC	mg/kg	27	18.6	14.9	33.7
57-12-5	CYANIDE	mg/kg	0.042 J	0.434	0.069 J	0.093 J

		Dup of SB-14(45-47)						
Con Ed - Ludl	ow Site	Location ID:	SB-14	SB-14	SB-14	SB-16		
Validated Soil	Analytical Data	Sample ID:	SB-14(38-40)	SB-14(45-47)	SB-14A(45-47)	SB-16(30-33)		
SDGs: D4710	, D1751, D5300	Lab Sample Id:	D5300-03	D5300-06	D5300-07	D5300-01		
		Depth:	38 - 40 ft	45 - 47 ft	45 - 47 ft	30 - 33 ft		
		Source:	CTECH	CTECH	CTECH	CTECH		
		SDG:	D5300	D5300	D5300	D5300		
		Matrix:	SOIL	SOIL	SOIL	SOIL		
		Sampled:	12/21/2012 10:50	12/21/2012 11:50	12/21/2012 12:00	12/20/2012 16:10		
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013		
CAS NO.	COMPOUND	UNITS:						
	VOLATILES							
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	1.1 U	1 U	1 U	110 U		
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.59 U	0.54 U	0.53 U	57.5 U		
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	1.7 U	1.5 U	1.5 U	170 U		
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.2 U	1 U	1 U	110 U		
75-34-3	1,1-DICHLOROETHANE	ug/kg	1.2 U	1.1 U	1.1 U	120 U		
75-35-4	1,1-DICHLOROETHENE	ug/kg	1.9 U	1.7 U	1.7 U	180 U		
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.64 UJ	0.58 U	0.58 U	62.5 U		
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.9 U	0.82 UJ	0.81 UJ	87.5 U		
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1.1 U	R	R	110 U		
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/kg	0.82 U	0.75 U	0.74 U	80 U		
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.79 U	0.72 U	0.72 U	77.5 U		
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.82 U	0.75 U	0.74 U	80 U		
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.33 U	0.3 U	0.3 U	32.5 U		
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.47 U	0.43 U	0.43 U	46.3 U		
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.52 U	0.48 U	0.47 U	51.3 U		
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	64 U	58 UJ	58 UJ	1300 UJ		
591-78-6	2-HEXANONE	ug/kg	5 U	4.6 U	4.5 U	490 UJ		
67-64-1	ACETONE	ug/kg	3.9 U	27 J	26 J	380 U		
71-43-2	BENZENE	ug/kg	0.49 U	0.44 U	0.44 U	17200		
74-97-5	BROMOCHLOROMETHANE	ug/kg	1 U	0.92 U	0.91 U	98.8 U		
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.79 U	0.72 U	0.72 U	77.5 U		
75-25-2	BROMOFORM	ug/kg	0.95 U	0.86 U	0.85 U	92.5 UJ		
74-83-9	BROMOMETHANE	ug/kg	3.1 UJ	2.9 U	2.8 U	310 U		
75-15-0	CARBON DISULFIDE	ug/kg	1.4 U	1.2 U	1.2 U	130 U		
56-23-5	CARBON TETRACHLORIDE	ug/kg	1.3 U	1.2 U	1.1 U	120 U		
108-90-7	CHLOROBENZENE	ug/kg	0.64 U	0.58 U	0.58 U	62.5 U		
75-00-3	CHLOROETHANE	ug/kg	1.8 U	1.6 U	1.6 U	180 U		
67-66-3	CHLOROFORM	ug/kg	0.95 U	0.86 U	0.85 U	92.5 U		
74-87-3	CHLOROMETHANE	ug/kg	1.1 U	1 U	0.99 U	320 J		
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	1.1 U	1 U	1 U	110 U		
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.92 U	0.84 U	0.83 U	90 U		
110-82-7	CYCLOHEXANE	ug/kg	1.3 U	1.2 U	1.2 U	130 U		
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.69 U	0.63 UJ	0.62 UJ	67.5 UJ		
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.83 U	0.76 UJ	0.75 UJ	81.3 U		
100-41-4	ETHYLBENZENE	ug/kg	0.79 U	0.72 U	0.72 U	43900		
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.61 U	0.56 U	0.55 U	3300		
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.92 U	0.84 U	0.83 U	38400		
79-20-9	METHYL ACETATE	ug/kg	1.9 U	1.8 U	1.7 U	190 U		
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/kg	4 U	3.6 U	3.6 U	390 UJ		
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/kg	3.7 U	3.4 U	3.4 U	370 U		
108-87-2	METHYLCYCLOHEXANE	ug/kg	1.4 U	1.2 U	1.2 U	180 J		
75-09-2	METHYLENE CHLORIDE	ug/kg	1.8 UJ	1.7 U	1.6 U	180 U		
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.87 U	0.79 U	0.78 U	18200		
100-42-5	STYRENE	ug/kg	0.58 U	0.52 U	0.52 U	8100		
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	1.2 U	1.1 U	1.1 U	120 U		
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	1.3 U	1.2 U	1.2 U	130 U		
108-88-3	TOLUENE	ug/kg	0.82 U	0.75 U	0.74 U	46500		
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.88 U	0.8 U	0.8 U	86.3 U		
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	1 U	0.92 U	0.91 U	98.8 U		
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	1.1 U	1 U	0.99 U	110 U		
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg	1.7 U	1.5 U	1.5 U	170 U		
75-01-4	VINYL CHLORIDE	ug/kg	1.6 U	1.4 U	1.4 U	150 U		

					Dup of SB-14(45-47)	
Con Ed - Ludlo	w Site	Location ID:	SB-14	SB-14	SB-14	SB-16
Validated Soil	Analytical Data	Sample ID:	SB-14(38-40)	SB-14(45-47)	SB-14A(45-47)	SB-16(30-33)
SDGs: D4710,	D1751, D5300	Lab Sample Id:	D5300-03	D5300-06	D5300-07	D5300-01
		Depth:	38 - 40 ft	45 - 47 ft	45 - 47 ft	30 - 33 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D5300	D5300	D5300	D5300
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled: Validated:	12/21/2012 10:50 1/23/2013	12/21/2012 11:50 1/23/2013	12/21/2012 12:00 1/23/2013	12/20/2012 16:10 1/23/2013
CAS NO.	COMPOUND	UNITS:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
	SEMIVOLATILES	UNIIS.				
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	16.7 U	15.2 U	15.1 U	16.4 U
	2,3,4,6-TETRACHLOROPHENOL	ug/kg	16.7 U	15.2 U	15.1 U	16.4 U
	2,4,5-TRICHLOROPHENOL	ug/kg	29.9 U	27.1 U	26.9 U	29.2 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	13 U	11.8 U	11.7 U	12.7 U
	2,4-DICHLOROPHENOL	ug/kg	16.2 U	14.7 U	14.6 U	15.9 U
	2,4-DIMETHYLPHENOL	ug/kg	24.2 U	21.9 U	21.7 U	23.6 U
	2,4-DINITROPHENOL	ug/kg	43.3 UJ	39.3 UJ	39 UJ	42.3 UJ
	2,4-DINITROTOLUENE	ug/kg	12.9 U	11.7 U	11.6 U	12.6 U
	2,6-DINITROTOLUENE	ug/kg	17.4 U	15.8 U	15.6 U	17 U
	2-CHLORONAPHTHALENE	ug/kg	9.7 U	8.8 U	8.7 U	9.5 U 22 U
	2-CHLOROPHENOL 2-METHYLNAPHTHALENE	ug/kg ug/kg	22.5 U 10.7 U	20.4 U 9.7 U	20.3 U 9.7 U	14900
	2-METHYLPHENOL (O-CRESOL)	ug/kg ug/kg	23.1 U	21 U	20.8 U	22.6 U
	2-NITROANILINE	ug/kg ug/kg	18.9 U	17.1 U	17 U	18.5 U
	2-NITROPHENOL	ug/kg	20.6 U	18.6 U	18.5 U	20.1 U
	3,3'-DICHLOROBENZIDINE	ug/kg	27.3 U	24.8 U	24.6 U	26.7 U
МЕРН3МЕРН	3+4-Methylphenols	ug/kg	22.1 U	20 U	19.9 U	21.6 U
	3-NITROANILINE	ug/kg	27.3 U	24.8 U	24.6 U	26.7 U
	4,6-DINITRO-2-METHYLPHENOL	ug/kg	24.4 UJ	22.1 UJ	22 UJ	23.9 UJ
	4-BROMOPHENYL PHENYL ETHER	ug/kg	8.3 U	7.5 U	7.5 U	8.1 U
	4-CHLORO-3-METHYLPHENOL	ug/kg	18.9 U	17.1 U	17 U	18.5 U
	4-CHLOROANILINE	ug/kg	30 U	27.2 U	27 U	29.4 U
	4-CHLOROPHENYL PHENYL ETHER 4-NITROANILINE	ug/kg	23.1 U	21 U	20.8 U	22.6 U 54.2 U
	4-NITROANILINE 4-NITROPHENOL	ug/kg ug/kg	55.5 U 79.1 U	50.3 U 71.7 U	49.9 U 71.2 U	77.3 U
	ACENAPHTHENE	ug/kg ug/kg	12 U	10.9 U	10.8 U	1800
208-96-8	ACENAPHTHYLENE	ug/kg	10.7 U	9.7 U	9.7 U	4700
	ACETOPHENONE	ug/kg	13 U	11.8 U	11.7 U	12.7 U
	ANTHRACENE	ug/kg	8.7 U	7.9 U	7.8 U	2500
	ATRAZINE	ug/kg	22.5 U	20.4 U	20.3 U	22 U
	BENZALDEHYDE	ug/kg	22.2 U	20.2 U	20 U	21.7 U
	BENZO(A)ANTHRACENE	ug/kg	20.3 U	18.4 U	18.3 U	1600
	BENZO(A)PYRENE	ug/kg	9.2 U	8.3 U	8.3 U	1200
	BENZO(B)FLUORANTHENE	ug/kg	13.9 U	12.6 U	12.5 U	940
	BENZO(G,H,I)PERYLENE BENZO(K)FLUORANTHENE	ug/kg ug/kg	17.3 U 20.1 U	15.6 U 18.2 U	15.5 U 18.1 U	310 J 380 J
	BENZYL BUTYL PHTHALATE	ug/kg ug/kg	20.1 U 20.4 U	18.5 U	18.4 U	20 U
	BIPHENYL (DIPHENYL)	ug/kg	16.1 U	14.6 U	14.5 U	1600
	BIS(2-CHLOROETHOXY) METHANE	ug/kg	24.5 U	22.2 U	22.1 U	24 U
	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/kg	20.4 U	18.5 U	18.4 U	20 U
	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	17.6 U	16 U	15.9 U	17.2 U
	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	15.1 U	13.7 U	13.6 U	14.7 U
	CAPROLACTAM	ug/kg	19.8 U	18 U	17.8 U	19.4 U
	CARBAZOLE	ug/kg	9.3 U	8.5 U	8.4 U	9.1 U
	CHRYSENE DIBENIZA HI ANTHIR ACENIE	ug/kg	19.3 U	17.5 U	17.4 U	1600
	DIBENZ(A,H)ANTHRACENE DIBENZOFURAN	ug/kg ug/kg	12.3 U 16.6 U	11.1 U 15.1 U	11 U 15 U	12 U 530
	DIETHYL PHTHALATE	ug/kg ug/kg	6.6 U	6 U	6 U	6.5 U
131-11-3	DIMETHYL PHTHALATE	ug/kg ug/kg	570	320 J	340 J	250 J
	DI-N-BUTYL PHTHALATE	ug/kg	33.5 U	30.3 U	30.1 U	32.7 U
	DI-N-OCTYLPHTHALATE	ug/kg	4.9 U	4.4 U	4.4 U	4.7 U
	FLUORANTHENE	ug/kg	8.6 U	7.8 U	7.7 U	2500
	FLUORENE	ug/kg	16.1 U	14.6 U	14.5 U	4500
	HEXACHLOROBENZENE	ug/kg	17.4 U	15.8 U	15.6 U	17 U
	HEXACHLOROBUTADIENE	ug/kg	15.5 U	14 U	13.9 U	15.1 U
	HEXACHLOROCYCLOPENTADIENE	ug/kg	10.4 UJ	9.4 UJ	9.3 UJ	10.1 UJ
	HEXACHLOROETHANE	ug/kg	19 U	17.3 U	17.1 U	18.6 U
	INDENO(1,2,3-C,D)PYRENE ISOPHORONE	ug/kg	14.2 U	12.9 U	12.8 U	200 J
	NAPHTHALENE	ug/kg ug/kg	14.1 U 14.7 U	12.7 U 13.3 U	12.7 U 13.2 U	13.7 U 33700
	NITROBENZENE	ug/kg ug/kg	16.1 U	14.6 U	13.2 U 14.5 U	15.7 U
	N-NITROSODI-N-PROPYLAMINE	ug/kg ug/kg	21.5 U	19.5 U	19.3 U	21 U
	N-NITROSODIPHENYLAMINE	ug/kg	10.2 U	9.3 U	9.2 U	10 U
	PENTACHLOROPHENOL	ug/kg	29.1 U	26.4 U	26.2 U	28.5 U
85-01-8	PHENANTHRENE	ug/kg	11.5 U	10.4 U	10.4 U	9900
	PHENOL	ug/kg	9.8 U	8.9 U	8.9 U	9.6 U
129-00-0	PYRENE	ug/kg	10.2 U	9.3 U	9.2 U	3300

					Dup of SB-14(45-47)	
Con Ed - Ludlow Site		Location ID:	SB-14	SB-14	SB-14	SB-16
Validated Soil Analytical Data		Sample ID:	SB-14(38-40)	SB-14(45-47)	SB-14A(45-47)	SB-16(30-33)
SDGs: D4710	, D1751, D5300	Lab Sample Id:	D5300-03	D5300-06	D5300-07	D5300-01
		Depth:	38 - 40 ft	45 - 47 ft	45 - 47 ft	30 - 33 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D5300	D5300	D5300	D5300
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	12/21/2012 10:50	12/21/2012 11:50	12/21/2012 12:00	12/20/2012 16:10
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	7400 J	2940 J	3060 J	14400 J
7440-36-0	ANTIMONY	mg/kg	0.59 UJ	0.53 UJ	0.55 UJ	0.59 UJ
7440-38-2	ARSENIC	mg/kg	0.97 J	0.71 J	0.89 J	5.98
7440-39-3	BARIUM	mg/kg	54.1 J	23.7 Ј	26.6 J	83.1 J
7440-41-7	BERYLLIUM	mg/kg	0.39	0.14 J	0.14 J	0.32 J
7440-43-9	CADMIUM	mg/kg	0.06 U	0.06 U	0.06 U	0.06 U
7440-70-2	CALCIUM	mg/kg	2180 J	9830 J	9770 J	24700 J
7440-47-3	CHROMIUM, TOTAL	mg/kg	16.3 J	7.26 UJ	7.44 UJ	12.4 J
7440-48-4	COBALT	mg/kg	7.86	3.26	3.53	3.36
7440-50-8	COPPER	mg/kg	7.99	5.12	5.88	5.21
7439-89-6	IRON	mg/kg	18000 J	8120 J	8770 J	15400 J
7439-92-1	LEAD	mg/kg	10.1	4.47	4.63	25.8
7439-95-4	MAGNESIUM	mg/kg	3730 J	4070 J	3950 J	13500 J
7439-96-5	MANGANESE	mg/kg	424 J	290 J	296 J	652 J
7439-97-6	MERCURY	mg/kg	0.005 J	0.002 U	0.003 J	0.003 J
7440-02-0	NICKEL	mg/kg	18.7	8.25	8.93	10.1
7440-09-7	POTASSIUM	mg/kg	2720 J	832 J	824 J	1040 J
7782-49-2	SELENIUM	mg/kg	0.76 J	0.39 U	0.41 U	0.49 J
7440-22-4	SILVER	mg/kg	0.51 J	0.17 J	0.15 J	0.23 J
7440-23-5	SODIUM	mg/kg	332	101	97.1 J	6400
7440-28-0	THALLIUM	mg/kg	0.28 U	0.25 U	0.27 U	0.29 U
7440-62-2	VANADIUM	mg/kg	19.3	8.87	10.2	19.1
7440-66-6	ZINC	mg/kg	26.9 J	13 J	12.9 J	12.6 J
57-12-5	CYANIDE	mg/kg	0.241 J	0.219 J	0.299	0.909

Con Ed - Ludl	ow Site	Location ID:	SB-16	SB-17	SB-17	TP-6FLOOR
Validated Soil Analytical Data		Sample ID:	SB-16(45-47)	SB-17(33-35)	SB-17(45-47)	TP-6FLOOR
SDGs: D4710, D1751, D5300		Lab Sample Id:	D5300-02	D5300-08	D5300-09	D4751-10
5D G3. D 4710.	, 51751, 55500	Depth:	45 - 47 ft	33 - 35 ft	45 - 47 ft	D-1/31 10 -
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D5300	D5300	D5300	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	12/20/2012 17:10	12/21/2012 15:00	12/21/2012 15:30	11/7/2012 10:30
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:		.,,,		
CLIDITO.	VOLATILES	CITIE.				
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	1.1 U	0.92 U	0.95 U	1.3 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.58 U	0.48 U	0.5 U	R
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	1.7 U	1.4 U	1.4 U	1.9 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.1 U	0.94 U	0.98 U	1.3 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	1.2 U	0.98 U	1 U	6.6 J
75-35-4	1,1-DICHLOROETHENE	ug/kg	1.9 U	1.5 U	1.6 U	2.1 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.63 U	0.52 U	0.54 UJ	R
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.88 UJ	0.73 UJ	0.76 U	R
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	R	R	0.94 U	R
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/kg	0.81 U	0.67 U	0.69 U	0.92 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.78 U	0.65 U	0.67 U	R
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.81 U	0.67 U	0.69 U	0.92 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.33 U	0.27 U	0.28 U	0.37 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.47 U	0.38 U	0.4 U	R
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.52 U	0.43 U	0.44 U	R
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	63 UJ	10.4 UJ	54 U	70 U
591-78-6	2-HEXANONE	ug/kg	5 U	4.1 U	4.3 U	5.6 U
67-64-1	ACETONE	ug/kg	71	3.1 U	11 J	26 J
71-43-2	BENZENE	ug/kg	0.48 U	0.4 U	0.41 U	0.54 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	1 U	0.82 U	0.86 U	1.1 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.78 U	0.65 U	0.67 U	0.89 U
75-25-2	BROMOFORM	ug/kg	0.93 U	0.77 U	0.8 U	1.1 UJ
74-83-9	BROMOMETHANE	ug/kg	3.1 U	2.5 U	2.7 UJ	3.5 U
75-15-0	CARBON DISULFIDE	ug/kg	1.3 U	1.1 U	1.1 U	1.5 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	1.3 U	1 U	1.1 U	1.4 U
108-90-7	CHLOROBENZENE	ug/kg	0.63 U	0.52 U	0.54 U	0.72 UJ
75-00-3	CHLOROETHANE	ug/kg	1.8 U	1.5 U	1.5 U	2 U
67-66-3	CHLOROFORM	ug/kg	0.93 U	0.77 U	0.8 U	1.1 U
74-87-3	CHLOROMETHANE	ug/kg	1.1 U	0.89 U	0.93 U	1.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	1.1 U	0.93 U	0.97 U	1.3 U
10061-01-5 110-82-7	CIS-1,3-DICHLOROPROPENE CYCLOHEXANE	ug/kg	0.91 U	0.75 U	0.78 U	1 U
		ug/kg	1.3 U	1.1 U	1.1 U	1.4 U
124-48-1 75-71-8	DIBROMOCHLOROMETHANE DICHLORODIFLUOROMETHANE	ug/kg	0.68 UJ 0.82 UJ	0.56 UJ 0.68 UJ	0.59 U 0.71 U	0.77 U 0.93 U
100-41-4	ETHYLBENZENE	ug/kg ug/kg	0.82 UJ 4.5 J	0.68 UJ 0.65 U	0.71 U 0.67 U	0.93 U 0.89 UJ
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg ug/kg	4.5 J 0.61 U	0.65 U	0.67 U 0.52 U	0.89 UJ R
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg ug/kg	3.9 J	0.5 U	0.78 U	1 UJ
79-20-9	METHYL ACETATE	ug/kg ug/kg	1.9 U	1.6 U	1.6 U	2.2 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/kg ug/kg	3.9 U	3.2 U	3.4 U	4.5 U
108-10-1	METHYL ISOBUTYL KETONE (2-BC FANONE) METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/kg ug/kg	3.7 U	3.2 U	3.4 U	4.3 U
108-10-1	METHYLCYCLOHEXANE	ug/kg	1.3 U	1.1 U	1.1 U	1.5 U
75-09-2	METHYLENE CHLORIDE	ug/kg ug/kg	1.8 U	1.5 U	1.5 UJ	7.2 UJ
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg ug/kg	2.2 J	0.71 U	0.74 U	0.97 UJ
100-42-5	STYRENE	ug/kg	0.57 U	0.47 U	0.49 U	0.64 UJ
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	1.2 U	1 U	1 U	1.4 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	1.3 U	1.1 U	1.1 U	1.4 UJ
108-88-3	TOLUENE	ug/kg	0.81 U	0.67 U	0.69 U	0.92 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.87 U	0.72 U	0.75 U	0.99 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	1 U	0.82 U	0.86 U	1.1 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	1.1 U	0.89 U	0.93 U	1.2 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg	1.7 U	1.4 U	1.4 U	1.9 U
75-01-4	VINYL CHLORIDE	ug/kg	1.6 U	1.3 U	1.3 U	1.8 U

Con Ed. Ludl	ow Cita	Logotion ID:	CD 16	CD 17	CD 17	TD 6ELOOD
Con Ed - Ludle Validated Soil	ow Site Analytical Data	Location ID: Sample ID:	SB-16 SB-16(45-47)	SB-17 SB-17(33-35)	SB-17 SB-17(45-47)	TP-6FLOOR TP-6FLOOR
			D5300-02	D5300-08	D5300-09	D4751-10
DD G3. D +710,	, 51751, 55500	Lab Sample Id: Depth:	45 - 47 ft	33 - 35 ft	45 - 47 ft	D-1/31 10 -
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D5300	D5300	D5300	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	12/20/2012 17:10	12/21/2012 15:00	12/21/2012 15:30	11/7/2012 10:30
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	16.6 U	13.6 U	14.2 U	19 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	16.6 U	13.6 U	14.2 U	19 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	29.7 U	24.2 U	25.3 U	34 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	12.9 U	10.6 U	11 U	15 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	16.1 U	13.1 U	13.7 U	18 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	24 U	19.6 U	20.4 U	27 U
51-28-5 121-14-2	2,4-DINITROPHENOL 2,4-DINITROTOLUENE	ug/kg	43 UJ 12.8 U	35.1 UJ 10.5 U	36.6 UJ 10.9 U	49 UJ 15 U
606-20-2	2,6-DINITROTOLUENE	ug/kg ug/kg	17.2 U	14.1 U	14.7 U	20 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg ug/kg	9.6 U	7.9 U	8.2 U	11 U
95-57-8	2-CHLOROPHENOL	ug/kg	22.3 U	18.2 U	19 U	25 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	10.6 U	8.7 U	190 J	12 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	22.9 U	18.7 U	19.6 U	26 U
88-74-4	2-NITROANILINE	ug/kg	18.8 U	15.3 U	16 U	21 U
88-75-5	2-NITROPHENOL	ug/kg	20.4 U	16.7 U	17.4 U	23 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	27.1 U	22.2 U	23.1 U	31 U
	3+4-Methylphenols	ug/kg	21.9 U	17.9 U	18.7 U	25 U
99-09-2	3-NITROANILINE	ug/kg	27.1 U	22.2 U	23.1 U	31 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	24.2 UJ	19.8 UJ	20.6 UJ	28 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	8.2 U	6.7 U	7 U	9.4 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	18.8 U	15.3 U	16 U	21 U
106-47-8	4-CHLOROANILINE	ug/kg	29.8 U	24.3 U	25.4 U	34 UJ
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	22.9 U	18.7 U	19.6 U	26 U
100-01-6 100-02-7	4-NITROANILINE	ug/kg	55 U	44.9 U	46.9 U	63 U 90 U
83-32-9	4-NITROPHENOL ACENAPHTHENE	ug/kg ug/kg	78.4 U 11.9 U	64.1 U 9.7 U	66.9 U 10.2 U	90 U 14 U
208-96-8	ACENAPHTHYLENE	ug/kg ug/kg	10.6 U	8.7 U	9.1 U	14 U
98-86-2	ACETOPHENONE	ug/kg	12.9 U	10.6 U	11 U	15 U
120-12-7	ANTHRACENE	ug/kg	8.6 U	7 U	7.4 U	9.8 U
1912-24-9	ATRAZINE	ug/kg	22.3 U	18.2 U	19 U	25 U
100-52-7	BENZALDEHYDE	ug/kg	22.1 U	18 U	18.8 U	25 U
56-55-3	BENZO(A)ANTHRACENE	ug/kg	20.2 U	16.5 U	17.2 U	350 J
50-32-8	BENZO(A)PYRENE	ug/kg	9.1 U	7.5 U	7.8 U	410 J
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	13.8 U	11.3 U	11.8 U	500
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	17.1 U	14 U	14.6 U	290 J
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	19.9 U	16.3 U	17 U	23 U
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	20.3 U	16.6 U	17.3 U	23 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	16 U	13 U	13.6 U	18 U
111-91-1 111-44-4	BIS(2-CHLOROETHOXY) METHANE	ug/kg	24.3 U 20.3 U	19.9 U	20.8 U 17.3 U	28 U 23 U
108-60-1	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER) BIS(2-CHLOROISOPROPYL) ETHER	ug/kg ug/kg	20.3 U 17.5 U	16.6 U 14.3 U	17.3 U 14.9 U	23 U 20 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg ug/kg	17.5 U 15 U	14.3 U 12.2 U	480	20 U 17 U
105-60-2	CAPROLACTAM	ug/kg ug/kg	19.6 U	16 U	16.8 U	22 U
86-74-8	CARBAZOLE	ug/kg	9.3 U	7.6 U	7.9 U	11 U
218-01-9	CHRYSENE	ug/kg	19.1 U	15.6 U	16.3 U	400 J
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	12.2 U	9.9 U	10.4 U	14 U
132-64-9	DIBENZOFURAN	ug/kg	16.5 U	13.5 U	14.1 U	19 U
84-66-2	DIETHYL PHTHALATE	ug/kg	6.6 U	5.4 U	5.6 U	7.5 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	400 J	350	390	632
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	33.2 U	27.1 U	150 J	38 U
117-84-0	DI-N-OCTYLPHTHALATE	ug/kg	4.8 U	3.9 U	4.1 U	5.5 U
206-44-0	FLUORANTHENE	ug/kg	8.5 U	6.9 U	7.2 U	750
86-73-7	FLUORENE HEYACHI OROBENZENE	ug/kg	16 U	13 U	13.6 U	18 U
118-74-1	HEXACHLOROBENZENE	ug/kg	17.2 U	14.1 U	14.7 U	20 U
87-68-3 77-47-4	HEXACHLOROBUTADIENE HEXACHLOROCYCLOPENTADIENE	ug/kg ug/kg	15.3 U 10.3 UJ	12.5 U 8.4 UJ	13.1 U 8.8 UJ	18 U 12 U
67-72-1	HEXACHLOROETHANE	ug/kg ug/kg	10.3 UJ 18.9 U	8.4 UJ 15.4 U	8.8 UJ 16.1 U	12 U 22 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg ug/kg	18.9 U 14.1 U	13.4 U 11.5 U	16.1 U	270 J
78-59-1	ISOPHORONE	ug/kg ug/kg	13.9 U	11.3 U 11.4 U	11.9 U	16 U
91-20-3	NAPHTHALENE	ug/kg ug/kg	14.6 U	11.4 U	250 J	10 U
98-95-3	NITROBENZENE	ug/kg	16 U	13 U	13.6 U	17 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	21.3 U	17.4 U	18.2 U	24 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	10.1 U	8.3 U	8.6 U	12 U
87-86-5	PENTACHLOROPHENOL	ug/kg	28.9 U	23.6 U	24.6 U	33 U
85-01-8	PHENANTHRENE	ug/kg	11.4 U	9.3 U	9.7 U	450 J
108-95-2	PHENOL	ug/kg	9.8 U	8 U	8.3 U	11 U
129-00-0	PYRENE	ug/kg	10.1 U	8.3 U	8.6 U	690

Con Ed - Luc	llow Site	Location ID:	SB-16	SB-17	SB-17	TP-6FLOOR
Validated So	il Analytical Data	Sample ID:	SB-16(45-47)	SB-17(33-35)	SB-17(45-47)	TP-6FLOOR
SDGs: D471	0, D1751, D5300	Lab Sample Id:	D5300-02	D5300-08	D5300-09	D4751-10
		Depth:	45 - 47 ft	33 - 35 ft	45 - 47 ft	-
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D5300	D5300	D5300	D4751
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	12/20/2012 17:10	12/21/2012 15:00	12/21/2012 15:30	11/7/2012 10:30
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	6040 J	2860 J	9120 J	9100 J
7440-36-0	ANTIMONY	mg/kg	0.61 UJ	0.47 UJ	0.51 UJ	0.36 UJ
7440-38-2	ARSENIC	mg/kg	1.18	0.68 J	0.73 J	14.5 J
7440-39-3	BARIUM	mg/kg	74.4 J	33.3 J	50.8 J	122 J
7440-41-7	BERYLLIUM	mg/kg	0.33	0.1 J	0.06 J	0.04 U
7440-43-9	CADMIUM	mg/kg	0.07 U	0.05 U	0.05 U	1.73
7440-70-2	CALCIUM	mg/kg	12500 J	1980 J	5560 J	6580 J
7440-47-3	CHROMIUM, TOTAL	mg/kg	12.9 J	10.6 UJ	25.5 J	28.7 J
7440-48-4	COBALT	mg/kg	7.11	4.01	6.27	9.95
7440-50-8	COPPER	mg/kg	9.59	8.81	31.1	173 J
7439-89-6	IRON	mg/kg	15800 J	7530 J	13000 J	39400
7439-92-1	LEAD	mg/kg	8.35	2.92	20.3	356
7439-95-4	MAGNESIUM	mg/kg	5840 J	2040 J	3950 J	4080 J
7439-96-5	MANGANESE	mg/kg	443 J	519 J	332 J	345 J
7439-97-6	MERCURY	mg/kg	0.005 J	0.003 J	0.014	6.2
7440-02-0	NICKEL	mg/kg	19.1	11.8	17.6	29.2
7440-09-7	POTASSIUM	mg/kg	2050 J	730 J	1420 J	901 J
7782-49-2	SELENIUM	mg/kg	0.63 J	0.38 J	0.51 J	2.18
7440-22-4	SILVER	mg/kg	0.37 J	0.17 J	0.18 J	0.39
7440-23-5	SODIUM	mg/kg	1100	188	913	2090 J
7440-28-0	THALLIUM	mg/kg	0.29 U	0.23 U	0.25 U	1.46
7440-62-2	VANADIUM	mg/kg	16.6	10.3	22.4	26 J
7440-66-6	ZINC	mg/kg	25.1 J	13.1 J	28.6 J	315
57-12-5	CYANIDE	mg/kg	0.294 J	0.172 J	0.257 J	0.297 J

Con Ed - Ludle	ow Site	Location ID:	FIELDQC
	Analytical Data	Sample ID:	FB122112
	D1751, D5300	Lab Sample Id:	D5300-10
DD 00. D 1710,	21701, 20000	Depth:	-
		Source:	CTECH
		SDG:	D5300
		Matrix:	SOIL
		Sampled:	12/21/2012 15:30
		Validated:	1/23/2013
CAS NO.	COMPOUND	UNITS:	ug/L
	VOLATILES		
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	0.38 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.36 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.2 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	0.46 U
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/kg	0.41 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.45 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.48 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.32 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	10 UJ
591-78-6	2-HEXANONE	ug/kg	1.9 U
67-64-1	ACETONE BENZENE	ug/kg	0.5 U 0.32 U
71-43-2 74-97-5	BROMOCHLOROMETHANE	ug/kg	0.32 U 0.2 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.2 U 0.36 U
75-27-4	BROMOFORM	ug/kg ug/kg	0.36 U 0.47 U
74-83-9	BROMOMETHANE	ug/kg ug/kg	0.47 U 0.2 UJ
75-15-0	CARBON DISULFIDE	ug/kg ug/kg	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/kg ug/kg	0.2 UJ
108-90-7	CHLOROBENZENE	ug/kg ug/kg	0.49 U
75-00-3	CHLOROETHANE	ug/kg ug/kg	0.2 U
67-66-3	CHLOROFORM	ug/kg	0.34 U
74-87-3	CHLOROMETHANE	ug/kg	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.31 U
110-82-7	CYCLOHEXANE	ug/kg	0.2 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.2 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.2 U
100-41-4	ETHYLBENZENE	ug/kg	0.2 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.45 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.95 U
79-20-9	METHYL ACETATE	ug/kg	0.2 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/kg	1.3 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/kg	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.2 U
75-09-2	METHYLENE CHLORIDE	ug/kg	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.43 U
100-42-5	STYRENE	ug/kg	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.35 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.27 U
108-88-3	TOLUENE TRANS 1.2 DICH OROSTHENE	ug/kg	0.37 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.41 U
10061-02-6 79-01-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.29 U 0.28 U
79-01-6 75-69-4	TRICHLOROETHYLENE (TCE) TRICHLOROFLUOROMETHANE	ug/kg ug/kg	0.28 U 0.35 U
75-09-4	VINYL CHLORIDE	ug/kg ug/kg	0.34 U
75 U1-T	THIL CILCONDE	ug/ng	0.57 0

Validance Sail Analytical Dana Sample ID: Lab Sample ID: Cab S	C E4 I41-	C'4-	IIti ID.	EIEI DOC
Display			Location ID:	FIELDQC
Depth: Source: CTECH Source: CTECH Source: CTECH D5300 Matrix: CTECH D5300				
CAS NO. COMPOUND	3DGs. D4710,	D1751, D5500		D3300-10
SDC:				CTECH
CAS NO. COMPOUND				
CAS NO. COMPOUND				
CAS NO. COMPOUND				12/21/2012 15:30
SEMINOLATILES				1/23/2013
15.94-13	CAS NO.	COMPOUND		ug/L
SS-90.2 2.3.4.6-TERACHLOROPHENOL ug/kg 0.48 U		SEMIVOLATILES		
95-95-4	95-94-3		ug/kg	
Section				
120.83-2 2.4-DICHLOROPHENOL				
105-67-9				
151-28-5				
121-14-2 2.4-DINTROTOLUENE		/		
506-20-2 2.6-DINTROTOLUENE Ug/Ng 0.38 U Ug/Ng 0.55-73 2.CHL.OROPHENOL Ug/Ng 0.64 U Ug/Ng 0.64 U Ug/Ng 0.64 U Ug/Ng 0.55-73 2.CHL.OROPHENOL Ug/Ng 0.58 U Ug/Ng 0.54 U Ug/Ng 0.54 U Ug/Ng 0.54 U Ug/Ng 0.54 U Ug/Ng 0.55 U Ug/				
91-58-7 2-CHLORONAPHTHALENE ug/kg				
95-57-8 2-CHLOROPHENOL				
91-57-6 2-METHYLNAPHTHALENE 91-58-49-7 2-METHYLNAPHTHALENE 91-58-7-5-5 2-METHYLPHENOL (O-CRESOL) 91-94-1 33-DICHLOROBENZIDINE 91-94-1 33-DICHLOROBENZIDINE 91-94-1 33-DICHLOROBENZIDINE 91-94-1 33-DICHLOROBENZIDINE 91-94-1 33-DICHLOROBENZIDINE 91-95-1 34-S2-1 4-DINITRO-2-METHYLPHENOL 91-95-2 4-DINITRO-2-METHYLPHENOL 91-95-2 95-50-7 4-CHLORO-3-METHYLPHENOL 91-95-2 95-50-7 4-CHLORO-3-METHYLPHENOL 91-95-2 95-50-7 4-CHLORO-3-METHYLPHENOL 91-95-2 91-96-47-8 4-CHLORO-3-METHYLPHENOL 91-95-2 91-90-1 91-00-1-6 4-NITRO-ANILINE 91-95-2 91-90-1 91-91-1 91-1 91-1-1 91-1 91-1-1 91-1 91-1-1 9				
95-48-7 2-METHYLPHENOL (O-CRESOL)	91-57-6			
S8-74-4 2-NITROANILINE	95-48-7			
18-75-5 2-NITROPHENOL ug/kg 0.62 U ug/kg 0.45 U ug/kg 0.88 U ug/kg 0.88 U ug/kg 0.88 U ug/kg 0.45 U ug/kg 0.48 U ug/kg 0.45 U ug/kg 0.48 U ug/kg 0.45 U ug/kg 0.45 U ug/kg 0.45 U ug/kg 0.45 U ug/kg 0.25 U ug/kg 0.48 U ug/kg 0.48 U ug/kg 0.49 U ug/kg 0.35 U ug/kg 0.25 U ug/kg 0.35 U ug/kg 0.30				
MEPHSMEPH 3-4-Methylphenols ug/kg 1.3 U ug/kg 1.4 ENDMOPHENYL PHENYL ETHER ug/kg 0.27 U ug/kg 1.4 CHILOROANILINE ug/kg 3.4 U	88-75-5	2-NITROPHENOL		0.62 U
99-09-2 3-NITROÁNLINE	91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	2.4 U
33-452-1 46-DINTRO-2-METHYLPHENOL ug/kg 0.27 U ug/kg 0.48 U ug/kg 0.49 U ug/kg 0.49 U ug/kg 0.25 U ug/kg 0.35 U				
101-55-3 4-BROMOPHENYL PHENYL ETHER ug/kg 0.48 U ug/kg 0.25 U ug/kg 0.26 U ug/kg 0.27 U ug/kg 0.28 U ug/kg 0.29 U ug/kg 0.29 U ug/kg 0.29 U ug/kg 0.29 U ug/kg 0.25 U	99-09-2			
59-50-7 4-CHLORO-3-METHYLPHENOL ug/kg ug/kg ug/kg 3.4 U ug/kg ug/kg 3.4 U ug/kg ug/kg 0.25 U ug/kg 0.19 U ug/kg 0.15 U ug/kg 0.25 U ug/kg				
106-47-8				
100-01-6 4-NITROANILINE				
100-01-6				
100-02-7				
SEA32-9 ACENAPHTHENE Ug/kg 0.25 U 208-96-8 ACENAPHTHYLENE Ug/kg 0.33 U Ug/kg 0.17 U Ug/kg 0.17 U Ug/kg 0.17 U Ug/kg 0.17 U Ug/kg 0.19 U Ug/kg 0.17 U Ug/kg 0.15 U Ug/kg 0.35 U Ug/kg 0.21 U Ug/kg 0.22 U Ug/kg 0.25 U Ug				
208-96-8 ACENAPHTHYLENE Ug/kg 0.83 U Ug/kg 0.17 U Ug/kg 0.19 U Ug/kg 0.92 U Ug/kg 0.95 U Ug/kg 0.35 U Ug/kg 0.21 U Ug/kg 0.23 U Ug/kg 0.25 U Ug/kg 0.65 U Ug/kg 0.				
98-86-2 ACETOPHENONE 120-12-7 ANTHRACENE 120-12-7 ANTHRACENE 1912-24-9 1912-24-9 1912-24-9 1912-24-9 1912-24-9 1912-24-9 1912-24-9 1912-24-2 1912-				
120-12-7 ANTHRACENE ug/kg 0.19 U 1912-24-9 ATRAZINE ug/kg 0.92 U 1912-24-9 ATRAZINE ug/kg 0.92 U 1912-24-9 BENZO(A)ANTHRACENE ug/kg 0.92 U 195-55-3 BENZO(A)ANTHRACENE ug/kg 0.17 U 205-99-2 BENZO(B)FLUORANTHENE ug/kg 0.35 U 191-24-2 BENZO(G)H,IDPERYLENE ug/kg 0.35 U 191-24-2 BENZO(G,H,DPERYLENE ug/kg 0.21 U 285-68-7 BENZYL BUTYL PHTHALATE ug/kg 0.21 U 285-68-7 BENZYL BUTYL PHTHALATE ug/kg 0.23 U 29-52-4 BIPHENYL (DIPHENYL) ug/kg 0.65 U 111-91-1 BIS(2-CHLOROETHOXY) METHANE ug/kg 0.65 U 1108-60-1 BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER) ug/kg 0.65 U 108-60-1 BIS(2-CHLOROISOPROPYL) ETHER ug/kg 0.19 U 105-60-2 CAPROLACTAM ug/kg 0.24 U 105-60-2 CAPROLACTAM ug/kg 0.26 U 218-01-9 CHRYSENE ug/kg 0.26 U 218-01-9 CHRYSENE ug/kg 0.21 U 131-01-3 DIBENZ(A,H)ANTHRACENE ug/kg 0.29 U 131-11-3 DIMETHYL PHTHALATE ug/kg 0.26 U 218-4-62 DIETHYL PHTHALATE ug/kg 0.26 U 219-131-11-3 DIMETHYL PHTHALATE ug/kg 0.26 U 217-84-72 DI-N-DCTYLPHTHALATE ug/kg 0.26 U 218-73-7 HUORANTHENE ug/kg 0.31 U 206-44-0 FLUORANTHENE ug/kg 0.31 U 206-44-0 FLUORANTHENE ug/kg 0.31 U 206-44-0 HEXACHLOROBUTADIENE ug/kg 0.31 U 206-44-0 HEXACHLOROPTENENE ug/kg 0.31 U 208-85-3 NITROBENZENE ug/kg 0.30 U 208-95-3 NITROBENZENE ug/kg 0.31 U 208-95-3 NITROBENZENE ug/kg 0.31 U 208-95-3 NITROBENZENE ug/kg 0.31 U 208-95-2 PHENOL ug/kg 0.31 U 208-95-2 Ug/kg 0.31 U 208-95-2 Ug/kg 0.31 U 208-95-2 Ug/kg 0.31 U 208-				
1912-24-9 ATRAZINE	120-12-7	ANTHRACENE		
56-55-3 BENZO(A)ANTHRACENE ug/kg 0.19 U 50-32-8 BENZO(A)PYRENE ug/kg 0.17 U 205-99-2 BENZO(G,H,DPERYLENE ug/kg 0.35 U 191-24-2 BENZO(K,FLUORANTHENE ug/kg 0.21 U 207-08-9 BENZO(K,FLUORANTHENE ug/kg 0.21 U 85-68-7 BENZYL BUTYL PHTHALATE ug/kg 0.21 U 92-52-4 BIPHENYL (DIPHENYL) ug/kg 0.65 U 111-91-1 BIS(2-CHLOROETHOXY) METHANE ug/kg 0.65 U 111-94-4 BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER) ug/kg 0.65 U 108-60-1 BIS(2-CHLOROISOPROPYL) ETHER ug/kg 0.65 U 108-60-2 BIS(2-CHLOROETHYL) PHTHALATE ug/kg 0.2 U 105-60-2 CAPOLACTAM ug/kg 0.2 U 86-74-8 CARBAZOLE ug/kg 0.2 U 218-01-9 CHRYSENE ug/kg 0.5 U 33-70-3 DIBENZOFURAN ug/kg 0.5 U 48-66-2 DIETHYL PHTHALATE ug/kg 0	1912-24-9	ATRAZINE		0.48 U
50-32-8 BENZO(A)PYRENE	100-52-7	BENZALDEHYDE	ug/kg	0.92 U
205-99-2 BENZO(B)FLUORANTHENE ug/kg 0.35 U 191-24-2 BENZO(G,H,DPERYLENE ug/kg 0.35 U 191-24-2 BENZO(K,JEUORANTHENE ug/kg 0.21 U 205-08-9 BENZO(K,JELUORANTHENE ug/kg 0.21 U 205-24 BIPHENYL (DIPHENYL) ug/kg 0.18 U 111-91-1 BIS(2-CHLOROETHOXY) METHANE ug/kg 0.65 U 111-44-1 BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER) ug/kg 0.65 U 108-60-1 BIS(2-CHLOROISOPROPYL) ETHER ug/kg 0.20 U 117-81-7 BIS(2-ETHYLHEXYL) PHTHALATE ug/kg 0.29 U 108-60-2 CAPROLACTAM ug/kg 0.26 U 218-01-9 CHRYSENE ug/kg 0.21 U 213-01-9 DIBENZO(A,H)ANTHRACENE ug/kg 0.21 U 213-01-9 DIBENZOCOPURAN ug/kg 0.25 U 24-66-2 DIETHYL PHTHALATE ug/kg 0.29 U 24-66-2 DIETHYL PHTHALATE ug/kg 0.26 U 211-11-3 DIMETHYL PHTHALATE ug/kg 0.26 U 24-74-2 DI-N-BUTYL PHTHALATE ug/kg 0.26 U 24-74-2 DI-N-BUTYL PHTHALATE ug/kg 0.26 U 206-44-0 FLUORANTHENE ug/kg 0.48 U 206-44-0 FLUORANTHENE ug/kg 0.49 U 206-44-0 FLUORANTHENE ug/kg 0.31 U 207-72-1 HEXACHLOROBUTADIENE ug/kg 0.31 U 208-3 HEXACHLOROBUTADIENE ug/kg 0.30 U 208-3 NAPHTHALENE ug/kg 0.31 U 209-3 NAPHTHALENE ug/kg 0.36 U 209-3 NAPHTHALENE ug/kg 0.31 U 208-0-3 NAPHTHALENE ug/kg 0.31 U 208-0-6 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 208-0-6 N-NITROSODI-N-PROPYLAMINE ug/kg 0.21 U 208-0-6 PENTACHLOROPHENOL ug/kg 0.31 U 208-0-6 PHENANTHRENE ug/kg 0.31 U 208-0-2 Ug/kg 0.31 U 208-0-2 PHENOL ug/kg 0.31 U 208-0-2 PHENOL ug/kg 0.31 U 208-0-2 Ug/kg 0.	56-55-3	BENZO(A)ANTHRACENE	ug/kg	0.19 U
191-24-2 BENZO(G,H,DPERYLENE Ug/kg Ug/				
207-08-9 BENZO(K)FLUORANTHENE ug/kg 0.21 U				
S5-68-7 BENZYL BUTYL PHTHALATE Ug/kg 0.23 U 92-52-4 BIPHENYL (DIPHENYL) Ug/kg 0.18 U 111-91-1 BIS(2-CHLOROETHOXY) METHANE Ug/kg 0.65 U 111-44-4 BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER) Ug/kg 0.65 U 108-60-1 BIS(2-CHLOROISOPROPYL) ETHER Ug/kg 0.2 U 107-60-2 CAPROLACTAM Ug/kg 0.26 U 108-67-8 CARBAZOLE Ug/kg 0.26 U 218-01-9 CHRYSENE Ug/kg 0.21 U 213-01-9 CHRYSENE Ug/kg 0.21 U 132-64-9 DIBENZ(A,H)ANTHRACENE Ug/kg 0.29 U 32-64-9 DIBENZOFURAN Ug/kg 0.29 U 34-74-2 DI-N-BUTYL PHTHALATE Ug/kg 0.26 U 34-74-2 DI-N-BUTYL PHTHALATE Ug/kg 0.26 U 217-84-0 DI-N-OCTYLPHTHALATE Ug/kg 0.26 U 34-74-2 DI-N-BUTYL PHTHALATE Ug/kg 0.26 U 38-73-7 FUORENE Ug/kg 0.48 U 38-673-7 FUORENE Ug/kg 0.37 U 118-74-1 HEXACHLOROBENZENE Ug/kg 0.3 U 37-68-3 HEXACHLOROCYCLOPENTADIENE Ug/kg 0.3 U 37-67-72-1 HEXACHLOROSTHANE Ug/kg 0.3 U 91-30-3 NAPHTHALENE Ug/kg 0.3 U 91-20-3 NAPHTHALENE				
92-52-4 BIPHENYL (DIPHENYL) 111-91-1 BIS(2-CHLOROETHOXY) METHANE 111-44-4 BIS(2-CHLOROETHOXY) METHANE 111-44-4 BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER) 108-60-1 BIS(2-CHLOROISOPROPYL) ETHER 117-81-7 BIS(2-ETHYLHEXYL) PHTHALATE 105-60-2 CAPROLACTAM 105-60-2 CAPROLACTAM 105-60-2 CARBAZOLE 118-01-9 CHRYSENE 1132-64-9 DIBENZ(A,H)ANTHRACENE 1132-64-9 DIBENZOFURAN 1132-64-9 DIBENZOFURAN 1132-64-9 DIBENZOFURAN 1131-11-3 DIMETHYL PHTHALATE 1131-11-3 DIMETHYL PHTHALATE 1131-11-3 DIMETHYL PHTHALATE 1131-11-3 DIMETHYL PHTHALATE 117-84-0 DI-N-OCTYLPHTHALATE 117-84-0 DI-N-OCTYLPHTHALATE 118-74-1 HEXACHLOROBENZENE 118-74-1 HEXACHLOROBUTADIENE 118-74-1 HEXACHLOROBUTADIENE 1193-39-5 INDENO(1,2,3-C,D)PYRENE 1193-39-5 INDENO(1,2,3-C,D)PYRENE 1191-20-3 NAPHTHALENE 1198-95-2 PHENOL 10.5 U ug/kg 10.18 U ug/kg 10.29 U ug/kg 10.31 U ug/kg 10.31 U 108-95-2 PHENOL 10.8 Ug/kg 10.18 U 10.8 Ug/kg 10.18 U 10.8 Ug/kg 10.18 U 10.8 Ug/kg 10.18 U 10.8 Ug/kg 10.19 U 10.8 Ug/kg 10.10 U 10.8 U 10.9 U 10.8 U 10.9 U 10.				
111-91-1 BIS(2-CHLOROETHOXY) METHANE				
111-44-4 BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER) 108-60-1 BIS(2-CHLOROISOPROPYL) ETHER 108-60-1 BIS(2-CHLOROISOPROPYL) ETHER 105-60-2 CAPROLACTAM 108-60-2 CAPROLACTAM 108-67-2 CARBAZOLE 108-67-8 1				
108-60-1 BIS(2-CHLOROISOPROPYL) ETHER				
117-81-7 BIS(2-ETHYLHEXYL) PHTHALATE ug/kg 0.19 U 105-60-2 CAPROLACTAM ug/kg 0.26 U ug/kg 0.26 U ug/kg 0.21 U ug/kg 0.21 U ug/kg 0.21 U ug/kg 0.25 U ug/kg 0.21 U ug/kg 0.25 U ug/kg 0.29 U ug/kg 0.29 U ug/kg 0.29 U ug/kg 0.26 U ug/kg 0.29 U ug/kg 0.29 U ug/kg 0.45 U ug/kg 0.26 U ug/kg 0.37 U ug/kg 0.31 U ug/kg 0.35 U ug/kg 0.31 U ug/kg 0.31 U ug/kg 0.35 U ug/kg 0.31 U ug/kg 0.35 U ug/kg 0.31 U ug/kg 0.35 U ug/kg 0.35 U ug/kg 0.31 U ug/kg 0.35 U ug/kg 0.35 U ug/kg 0.31 U ug/kg 0.35 U ug/kg 0.35 U ug/kg 0.31 U ug/kg 0.35 U ug/kg 0.35 U ug/kg 0.31 U ug/kg 0.35 U ug/kg				
105-60-2 CAPROLACTAM				
86-74-8 CARBAZOLE ug/kg 0.26 U 218-01-9 CHRYSENE ug/kg 0.21 U 53-70-3 DIBENZCA,H)ANTHRACENE ug/kg 0.29 U 132-64-9 DIBENZOFURAN ug/kg 0.29 U 84-66-2 DIETHYL PHTHALATE ug/kg 0.26 U 131-11-3 DIMETHYL PHTHALATE ug/kg 0.26 U 84-74-2 DI-N-BUTYL PHTHALATE ug/kg 0.61 U 206-44-0 FLUORANTHENE ug/kg 0.48 U 86-73-7 FLUORENE ug/kg 0.21 U 87-68-3 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 47-4-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.3 U 47-7-1 HEXACHLOROETHANE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 78-59-1 ISOPHORONE ug/kg 0.36 U 19-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.24 U 621-64-7 N-NITROSODI-N-PR		,		
53-70-3 DIBENZ(A,H)ANTHRACENE ug/kg 0.5 U 132-64-9 DIBENZOFURAN ug/kg 0.29 U 84-66-2 DIETHYL PHTHALATE ug/kg 0.45 U 131-11-3 DIMETHYL PHTHALATE ug/kg 0.26 U 84-74-2 DI-N-BUTYL PHTHALATE ug/kg 0.61 U 117-84-0 DI-N-OCTYLPHTHALATE ug/kg 0.48 U 206-44-0 FLUORANTHENE ug/kg 0.37 U 118-74-1 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 98-95-1 ISOPHORONE ug/kg 0.18 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U			ug/kg	0.26 U
132-64-9 DIBENZOFURAN ug/kg 0.29 U				
84-66-2 DIETHYL PHTHALATE ug/kg 0.45 U 131-11-3 DIMETHYL PHTHALATE ug/kg 0.26 U 84-74-2 DI-N-BUTYL PHTHALATE ug/kg 2.4 U 117-84-0 DI-N-OCTYLPHTHALATE ug/kg 0.61 U 206-44-0 FLUORANTHENE ug/kg 0.48 U 86-73-7 FLUORENE ug/kg 0.21 U 118-74-1 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 67-72-1 HEXACHLOROETHANE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 98-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.81 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODI-N-PROPYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.31 U				
131-11-3 DIMETHYL PHTHALATE Ug/kg 0.26 U				
84-74-2 DI-N-BUTYL PHTHALATE ug/kg 2.4 U 117-84-0 DI-N-OCTYLPHTHALATE ug/kg 0.61 U 206-44-0 FLUORANTHENE ug/kg 0.48 U 86-73-7 FLUORENE ug/kg 0.21 U 118-74-1 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 8-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 0.31 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
117-84-0 DI-N-OCTYLPHTHALATE ug/kg 0.61 U 206-44-0 FLUORANTHENE ug/kg 0.48 U 86-73-7 FLUORENE ug/kg 0.37 U 118-74-1 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 8-59-1 ISOPHORONE ug/kg 0.18 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 0.31 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
206-44-0 FLUORANTHENE ug/kg 0.48 U 86-73-7 FLUORENE ug/kg 0.37 U 118-74-1 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.29 U 67-72-1 HEXACHLOROETHANE ug/kg 0.18 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 8-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 0.31 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
86-73-7 FLUORENE ug/kg 0.37 U 118-74-1 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.29 U 67-72-1 HEXACHLOROETHANE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 78-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 0.31 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
118-74-1 HEXACHLOROBENZENE ug/kg 0.21 U 87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.29 U 67-72-1 HEXACHLOROETHANE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 78-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 0.21 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
87-68-3 HEXACHLOROBUTADIENE ug/kg 0.3 U 77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.29 U 67-72-1 HEXACHLOROETHANE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 78-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 2 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U	118-74-1		~ ~	
77-47-4 HEXACHLOROCYCLOPENTADIENE ug/kg 0.29 U 67-72-1 HEXACHLOROETHANE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 78-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 0.31 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U	87-68-3			
67-72-1 HEXACHLOROETHANE ug/kg 0.3 U 193-39-5 INDENO(1,2,3-C,D)PYRENE ug/kg 0.18 U 78-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 2 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U	77-47-4			
78-59-1 ISOPHORONE ug/kg 0.36 U 91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 2 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U	67-72-1	HEXACHLOROETHANE		0.3 U
91-20-3 NAPHTHALENE ug/kg 0.14 U 98-95-3 NITROBENZENE ug/kg 0.81 U 0621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 087-86-5 PENTACHLOROPHENOL ug/kg 0.71 U 087-86-5 PENTACHLOROPHENOL ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.31 U 089-95-2 Ug/kg 0.25 U			ug/kg	
98-95-3 NITROBENZENE ug/kg 0.81 U 02-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 08-30-6 N-NITROSODI-PROPYLAMINE ug/kg 0.71 U 08-78-86-5 PENTACHLOROPHENOL ug/kg 2 U 08-501-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U	78-59-1			
621-64-7 N-NITROSODI-N-PROPYLAMINE ug/kg 0.24 U 86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 2 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U	91-20-3			
86-30-6 N-NITROSODIPHENYLAMINE ug/kg 0.71 U 87-86-5 PENTACHLOROPHENOL ug/kg 2 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
87-86-5 PENTACHLOROPHENOL ug/kg 2 U 85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
85-01-8 PHENANTHRENE ug/kg 0.31 U 108-95-2 PHENOL ug/kg 0.25 U				
108-95-2 PHENOL ug/kg 0.25 U				
	129-00-0	PYRENE	ug/kg ug/kg	0.23 U 0.24 U

Con Ed - Ludl	ow Site	Location ID:	FIELDQC
Validated Soil	Analytical Data	Sample ID:	FB122112
SDGs: D4710,	D1751, D5300	Lab Sample Id:	D5300-10
		Depth:	-
		Source:	CTECH
		SDG:	D5300
		Matrix:	SOIL
		Sampled:	12/21/2012 15:30
		Validated:	1/23/2013
CAS NO.	COMPOUND	UNITS:	ug/L
	INORGANICS		
7429-90-5	ALUMINUM	mg/kg	44.8 J
7440-36-0	ANTIMONY	mg/kg	8 U
7440-38-2	ARSENIC	mg/kg	4.2 U
7440-39-3	BARIUM	mg/kg	4 U
7440-41-7	BERYLLIUM	mg/kg	0.7 U
7440-43-9	CADMIUM	mg/kg	0.5 U
7440-70-2	CALCIUM	mg/kg	599 J
7440-47-3	CHROMIUM, TOTAL	mg/kg	21.3
7440-48-4	COBALT	mg/kg	5.8 U
7440-50-8	COPPER	mg/kg	2 U
7439-89-6	IRON	mg/kg	551
7439-92-1	LEAD	mg/kg	2.6 U
7439-95-4	MAGNESIUM	mg/kg	35.6 J
7439-96-5	MANGANESE	mg/kg	5.61 J
7439-97-6	MERCURY	mg/kg	0.092 U
7440-02-0	NICKEL	mg/kg	7.24 J
7440-09-7	POTASSIUM	mg/kg	1190
7782-49-2	SELENIUM	mg/kg	4.8 U
7440-22-4	SILVER	mg/kg	1.5 U
7440-23-5	SODIUM	mg/kg	787 J
7440-28-0	THALLIUM	mg/kg	2.4 U
7440-62-2	VANADIUM	mg/kg	6.1 U
7440-66-6	ZINC	mg/kg	8.24 J
57-12-5	CYANIDE	mg/kg	3 U

ATTACHMENT A-2 VALIDATED LABORATORY DATA FOR GROUNDWATER SAMPLES

Con Ed - Ludlov			MW-1	MW-3	MW-4	MW-7	MW-8
Validated Group	ndwater Analytical Data	Location ID: Sample ID:	MW-1	MW-3	MW-4	MW-7	MW-8
vandated Groun	idwater Anaryticar Data	Lab Sample Id:	D4947-01	D4947-09	D4947-07	D4947-08	D4947-02
		Source:	CTECH	CTECH	CTECH	CTECH	CTECH
		SDG:	D4947	D4947	D4947	D4947	D4947
		Matrix:	WATER	WATER	WATER	WATER	WATER
		Sampled:	11/26/2012 10:30	11/27/2012 15:10	11/27/2012 11:50	11/27/2012 13:15	11/26/2012 13:20
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				1,20,200	
ľ	VOLATILES						
71-55-6	1,1,1-TRICHLOROETHANE	ug/l	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U	0.45 U
	1,1,2-TRICHLOROETHANE	ug/l	0.38 U	0.38 U	0.38 U	0.38 U	0.38 U
	1,1-DICHLOROETHANE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U	0.36 U
	1,1-DICHLOROETHENE	ug/l	0.47 U	0.47 U	0.47 U	0.47 U	0.47 U
	1,2,3-TRICHLOROBENZENE	ug/l	0.65 U	0.65 U	0.65 U	0.65 U	0.65 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/l	0.62 U	0.62 U	0.62 U	0.62 U	0.62 U
	1,2-DIBROMO-3-CHLOROPROPANE	ug/l	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/l	0.41 U	0.41 U	0.41 U	0.41 U	0.41 U
	1,2-DICHLOROBENZENE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U	0.45 U
	1,2-DICHLOROETHANE	ug/l	0.48 U	0.48 U	0.48 U	0.48 U	0.48 U
	1,2-DICHLOROPROPANE	ug/l	0.46 U	0.46 U	0.46 U	0.46 U	0.46 U
	1,3-DICHLOROBENZENE	ug/l	0.43 U	0.43 U	0.43 U	0.43 U	0.43 U
	1,4-DICHLOROBENZENE	ug/l	0.32 U	0.32 U	0.32 U	0.32 U	0.32 U
	1,4-DIOXANE (P-DIOXANE)	ug/l	50 UJ	50 UJ	50 U	50 UJ	50 UJ
	2-HEXANONE	ug/l	1.9 U	1.9 U	1.9 U	1.9 U	1.9 U
	ACETONE	ug/l	2.8 UJ	2.8 UJ	2.8 U	2.8 UJ	2.8 UJ
	BENZENE	ug/l	0.32 U	350	0.32 U	690	0.32 U
	BROMOCHLOROMETHANE	ug/l	2.2 U	2.2 U	2.2 U	2.2 U	2.2 U
	BROMODICHLOROMETHANE	ug/l	0.36 UJ	0.36 UJ	0.36 U	0.36 UJ	0.36 UJ
	BROMOFORM	ug/l	0.47 UJ	0.47 UJ	0.47 U	0.47 UJ	0.47 UJ
	BROMOMETHANE	ug/l	0.62 UJ	0.62 UJ	0.62 U	0.62 UJ	0.62 UJ
	CARBON DISULFIDE	ug/l	0.54 U	0.54 U	0.54 U	0.54 U	0.54 U
	CARBON TETRACHLORIDE	ug/l	0.62 UJ	0.62 UJ	0.62 U	0.62 UJ	0.62 UJ
	CHLOROBENZENE	ug/l	0.49 U	0.49 U	0.49 U	0.49 U	0.49 U
	CHLOROETHANE	ug/l	0.66 U	0.66 U	0.66 U	0.66 U	0.66 U
	CHLOROFORM	ug/l	0.34 U	0.34 U	0.34 U	0.34 U	0.34 U
	CHLOROMETHANE	ug/l	0.54 UJ	0.54 UJ	0.54 U	0.54 UJ	0.54 UJ
	CIS-1,2-DICHLOROETHYLENE	ug/l	0.35 U	0.45 J	0.35 U	0.35 U	0.35 U
	CIS-1,3-DICHLOROPROPENE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U	0.31 U
	CYCLOHEXANE	ug/l	0.55 U	5.5	0.55 U	2.7 J	0.55 U
	DIBROMOCHLOROMETHANE	ug/l	0.52 UJ	0.52 UJ	0.52 U	0.52 UJ	0.52 UJ
	DICHLORODIFLUOROMETHANE	ug/l	0.55 U	0.55 U	0.55 U	0.55 U	0.55 U
	ETHYLBENZENE (CHMENE)	ug/l	0.53 U	250	0.53 U	270	0.53 U 0.45 U
	ISOPROPYLBENZENE (CUMENE) M.P. YVI ENE (SUM OF ISOMEDS)	ug/l	0.45 U 0.95 U	7 410	0.45 U 0.95 U	24 420	0.45 U 0.95 U
	M,P-XYLENE (SUM OF ISOMERS) METHYL ACETATE	ug/l	0.95 U 0.83 UJ	0.83 UJ	0.95 U 0.83 U	0.83 UJ	0.95 U 0.83 UJ
	METHYL ACETATE METHYL ETHYL KETONE (2-BUTANONE)	ug/l	0.83 UJ 1.3 U	0.83 UJ 1.3 U	0.83 U 1.3 U	0.83 UJ 1.3 U	0.83 UJ 1.3 U
	METHYL ISOBUTYL KETONE (2-BUTANONE) METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/l ug/l	2.1 UJ	2.1 UJ	2.1 U	2.1 UJ	2.1 UJ
	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE) METHYLCYCLOHEXANE	ug/l ug/l	0.68 U	10	0.68 U	3.5 J	0.68 U
	METHYLENE CHLORIDE	_	0.41 U	0.41 U	0.41 U	0.41 U	0.68 U 0.41 U
	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/l ug/l	0.41 U 0.43 U	190	0.41 U 0.43 U	190	0.41 U 0.43 U
	STYRENE	ug/l ug/l	0.45 U	59	0.45 U	18	0.45 U
	TERT-BUTYL METHYL ETHER	ug/l ug/l	0.35 U	0.35 U	0.35 U	0.35 U	0.50 U 0.52 J
	TETRACHLOROETHYLENE(PCE)	ug/l ug/l	0.33 U 0.27 U	0.42 J	0.33 U 0.27 U	0.33 U 0.27 U	0.32 J 0.27 U
	TOLUENE	ug/l	0.27 U	280	0.27 U	270	0.27 U
	TRANS-1,2-DICHLOROETHENE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U	0.41 U
	TRANS-1,3-DICHLOROPROPENE	ug/l	0.41 U 0.29 UJ	0.41 U 0.29 UJ	0.41 U 0.29 U	0.41 U 0.29 UJ	0.41 U 0.29 UJ
	TRICHLOROETHYLENE (TCE)	ug/l	0.28 U	0.28 U	0.28 U	0.28 U	0.28 U
	TRICHLOROFLUOROMETHANE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U	0.35 U
	VINYL CHLORIDE	ug/l	0.34 U	0.34 U	0.34 U	0.34 U	0.34 U

Con Ed - Ludlo		Location ID:	MW-1	MW-3	MW-4	MW-7	MW-8
Validated Grou	ndwater Analytical Data	Sample ID:	MW-1	MW-3	MW-4	MW-7	MW-8
		Lab Sample Id:	D4947-01	D4947-09	D4947-07	D4947-08	D4947-02
		Source: SDG:	CTECH D4947	CTECH D4947	CTECH D4947	CTECH D4947	CTECH D4947
		Matrix:	WATER	WATER	WATER	WATER	WATER
		Sampled:	11/26/2012 10:30	11/27/2012 15:10		11/27/2012 13:15	11/26/2012 13:20
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:					
95-94-3	SEMIVOLATILES		0.22 11	0.21 111	0.22 11	0.21 11	0.21 11
95-94-3 58-90-2	1,2,4,5-TETRACHLOROBENZENE 2,3,4,6-TETRACHLOROPHENOL	ug/l ug/l	0.22 U 0.22 U	0.21 UJ 0.21 UJ	0.33 U 0.33 U	0.21 U 0.21 U	0.21 U 0.21 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/l	0.45 U	0.42 UJ	0.66 U	0.42 U	0.42 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/l	0.63 U	0.58 UJ	0.92 U	0.59 U	0.59 U
120-83-2	2,4-DICHLOROPHENOL	ug/l	0.74 U	0.69 U	1.1 U	0.69 U	0.69 U
105-67-9	2,4-DIMETHYLPHENOL	ug/l	0.8 U	0.74 U	1.2 U	0.75 U	0.75 U
51-28-5	2,4-DINITROPHENOL	ug/l	2.4 UJ	2.2 UJ	3.4 UJ	2.2 UJ	2.2 UJ
121-14-2	2,4-DINITROTOLUENE	ug/l	1.2 U	1.1 UJ	1.7 U	1.1 U	1.1 U
606-20-2	2,6-DINITROTOLUENE	ug/l	0.36 U	0.33 UJ	0.52 U	0.34 U	0.34 U
91-58-7 95-57-8	2-CHLORONAPHTHALENE	ug/l	0.18 U	0.17 UJ	0.26 U 0.89 U	0.17 U 0.57 U	0.17 U
93-37-8 91-57-6	2-CHLOROPHENOL 2-METHYLNAPHTHALENE	ug/l ug/l	0.61 U 0.36 U	0.56 U 320	0.89 U 0.52 U	130	0.57 U 0.34 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/l	0.27 U	0.25 U	0.32 U	0.25 U	0.25 U
88-74-4	2-NITROANILINE	ug/l	0.55 U	0.51 UJ	0.8 U	0.52 U	0.52 U
88-75-5	2-NITROPHENOL	ug/l	0.58 U	0.54 U	0.85 U	0.55 U	0.55 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/l	2.2 U	2.1 U	3.3 U	2.1 U	2.1 U
МЕРНЗМЕРН	3+4-Methylphenols	ug/l	0.43 U	0.4 U	0.62 U	0.4 U	0.4 U
99-09-2	3-NITROANILINE	ug/l	1.2 U	1.1 UJ	1.8 U	1.1 U	1.1 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/l	0.83 UJ	0.77 UJ	1.2 UJ	0.78 UJ	0.78 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/l	0.26 U	0.24 UJ	0.38 U	0.24 U	0.24 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/l	0.45 U	0.42 U	0.66 U	0.42 U	0.42 U
106-47-8 7005-72-3	4-CHLOROANILINE 4-CHLOROPHENYL PHENYL ETHER	ug/l	3.2 U 0.24 U	2.1 U 0.22 UJ	4.7 U 0.34 U	3 U 0.22 U	3 U 0.22 U
100-01-6	4-VITROANILINE	ug/l ug/l	1.5 U	0.22 UJ 1.4 UJ	2.2 U	1.4 U	1.4 U
100-02-7	4-NITROPHENOL	ug/l	2.2 U	2.1 UJ	3.3 U	2.1 U	2.1 U
83-32-9	ACENAPHTHENE	ug/l	0.24 U	17.1 J	0.34 U	12.1	0.22 U
208-96-8	ACENAPHTHYLENE	ug/l	0.79 U	84.2	1.1 U	7.9 J	0.74 U
98-86-2	ACETOPHENONE	ug/l	0.16 U	0.15 U	0.23 U	5.9 J	0.15 U
120-12-7	ANTHRACENE	ug/l	0.18 U	10.6 J	0.26 U	0.17 U	0.17 U
1912-24-9	ATRAZINE	ug/l	0.45 U	0.42 UJ	0.66 U	0.42 U	0.42 U
100-52-7	BENZALDEHYDE	ug/l	0.87 U	0.8 U	1.3 U	0.81 U	0.81 U
56-55-3 50-32-8	BENZO(A)ANTHRACENE BENZO(A)PYRENE	ug/l	0.18 U 0.16 U	0.17 U 0.15 U	0.26 U 0.23 U	0.17 U 0.15 U	0.17 U 0.15 U
205-99-2	BENZO(B)FLUORANTHENE	ug/l ug/l	0.10 U	0.13 U	0.23 U 0.48 U	0.13 U	0.13 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.33 U	0.3 U	0.48 U	0.31 U	0.31 U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.2 U	0.19 U	0.3 U	0.19 U	0.19 U
85-68-7	BENZYL BUTYL PHTHALATE	ug/l	0.21 U	0.2 U	0.31 U	0.2 U	0.2 U
92-52-4	BIPHENYL (DIPHENYL)	ug/l	0.17 U	25.1 J	0.25 U	7.3 J	0.16 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/l	0.62 U	0.57 U	0.9 U	0.58 U	0.58 U
111-44-4	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/l	0.62 U	0.57 U	0.9 U	0.58 U	0.58 U
108-60-1 117-81-7	BIS(2-CHLOROISOPROPYL) ETHER	ug/l	0.19 U	0.18 U	0.28 U	0.18 U	0.18 U
105-60-2	BIS(2-ETHYLHEXYL) PHTHALATE CAPROLACTAM	ug/l	0.18 U 2.2 U	0.17 U 2.1 U	15.4 J 3.3 U	0.17 U 2.1 U	0.17 U 2.1 U
86-74-8	CARBAZOLE	ug/l ug/l	0.25 U	6.2 J	0.36 U	0.23 U	0.23 U
218-01-9	CHRYSENE	ug/l	0.23 U	0.19 U	0.3 U	0.19 U	0.19 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.47 U	0.44 U	0.69 U	0.44 U	0.44 U
132-64-9	DIBENZOFURAN	ug/l	0.27 U	4.5 J	0.39 U	0.25 U	0.25 U
84-66-2	DIETHYL PHTHALATE	ug/l	0.43 U	0.4 UJ	0.62 U	0.4 U	0.4 U
131-11-3	DIMETHYL PHTHALATE	ug/l	0.25 U	0.23 UJ	0.36 U	0.23 U	0.23 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	2.2 U	2.1 UJ	3.3 U	2.1 U	2.1 U
117-84-0	DI-N-OCTYLPHTHALATE	ug/l	0.57 U	0.53 U	0.84 U	0.54 U	0.54 U
206-44-0 86-73-7	FLUORANTHENE FLUORENE	ug/l	0.45 U 0.35 U	7.6 J 34.7 J	0.66 U 0.51 U	0.42 U 9.9 J	0.42 U 0.33 U
86-73-7 118-74-1	HEXACHLOROBENZENE	ug/l ug/l	0.35 U 0.2 U	0.19 UJ	0.51 U 0.3 U	9.9 J 0.19 U	0.33 U 0.19 U
87-68-3	HEXACHLOROBUTADIENE	ug/l ug/l	0.28 U	0.19 UJ 0.26 U	0.41 U	0.19 U 0.26 U	0.19 U 0.26 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/l	0.27 U	0.25 UJ	0.41 U	0.25 U	0.25 U
67-72-1	HEXACHLOROETHANE	ug/l	0.28 U	0.26 U	0.41 U	0.26 U	0.26 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/l	0.17 U	0.16 U	0.25 U	0.16 U	0.16 U
78-59-1	ISOPHORONE	ug/l	0.34 U	0.31 U	0.49 U	0.32 U	0.32 U
91-20-3	NAPHTHALENE	ug/l	0.13 U	2500	0.2 U	1300	0.13 U
98-95-3	NITROBENZENE	ug/l	0.76 U	0.71 U	1.1 U	0.72 U	0.72 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/l	0.22 U	0.21 U	0.33 U	0.21 U	0.21 U
86-30-6 87-86-5	N-NITROSODIPHENYLAMINE	ug/l	0.67 U	0.63 UJ	0.98 U	0.63 U	0.63 U
	PENTACHLOROPHENOL PHENANTHRENE	ug/l ug/l	1.9 U 0.29 U	1.8 UJ 46.2 J	2.8 U 0.43 U	1.8 U 14.2	1.8 U 0.27 U
	PHENOL	ug/l ug/l	0.24 U	0.22 U	0.43 U 0.34 U	0.22 U	0.27 U 0.22 U
108-95-2			U.4T U	0.44 0	U.JT U	U.44 U	0.22

Con Ed - Lud	low Site	Location ID:	MW-1	MW-3	MW-4	MW-7	MW-8
Validated Gro	oundwater Analytical Data	Sample ID:	MW-1	MW-3	MW-4	MW-7	MW-8
		Lab Sample Id:	D4947-01	D4947-09	D4947-07	D4947-08	D4947-02
		Source:	CTECH	CTECH	CTECH	CTECH	CTECH
		SDG:	D4947	D4947	D4947	D4947	D4947
		Matrix:	WATER	WATER	WATER	WATER	WATER
		Sampled:	11/26/2012 10:30	11/27/2012 15:10	11/27/2012 11:50	11/27/2012 13:15	11/26/2012 13:20
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:	,				
	INORGANICS						
7429-90-5	ALUMINUM	ug/l	331 J	72.9 UJ	161000 J	268 J	1860 J
7440-36-0	ANTIMONY	ug/l	4.98 J	5.31 J	4 UJ	6.86 J	5.06 J
7440-38-2	ARSENIC	ug/l	6.1	2.45 J	206	2.1 U	4.85 J
7440-39-3	BARIUM	ug/l	159 J	106 J	2170 J	109 J	104 J
7440-41-7	BERYLLIUM	ug/l	0.35 UJ	0.35 UJ	7.1 J	0.35 UJ	0.35 UJ
7440-43-9	CADMIUM	ug/l	0.25 U	0.32 J	25	0.25 U	0.3 J
7440-70-2	CALCIUM	ug/l	91200 J	58900 J	67500 J	105000 J	78800 J
7440-47-3	CHROMIUM, TOTAL	ug/l	55.8 UJ	13.8 UJ	352 J	23.5 UJ	28.6 UJ
7440-48-4	COBALT	ug/l	2.9 U	2.9 U	263	2.9 U	4.32 J
7440-50-8	COPPER	ug/l	6.08 J	7.44 J	822 J	1.44 J	20.7 J
7439-89-6	IRON	ug/l	5890 J	1160 J	520000 J	680 UJ	3790 J
7439-92-1	LEAD	ug/l	1.34 J	6.72	10700	1.72 J	24.5 J
7439-95-4	MAGNESIUM	ug/l	22100 J	21500 J	53000 J	37800 J	25200 J
7439-96-5	MANGANESE	ug/l	4090 J	1100 J	6820 J	1030 J	1520 J
7439-97-6	MERCURY	ug/l	0.092 J	0.092 U	16.85	0.092 U	0.092 U
7440-02-0	NICKEL	ug/l	31.6 U	12.2 U	689	19.4 U	18.2 UJ
7440-09-7	POTASSIUM	ug/l	8980 J	8370 J	23300 J	8420 J	8280 J
7782-49-2	SELENIUM	ug/l	2.4 U	2.4 U	26	2.4 U	2.4 U
7440-22-4	SILVER	ug/l	0.75 UJ	0.75 UJ	25.7 J	0.75 UJ	0.75 UJ
7440-23-5	SODIUM	ug/l	361000 J	3480000 J	167000 J	535000 J	628000 J
7440-28-0	THALLIUM	ug/l	1.2 UJ	1.2 UJ	5.2 J	1.2 UJ	1.2 UJ
7440-62-2	VANADIUM	ug/l	3.05 UJ	3.05 UJ	438 J	3.05 UJ	5.36 J
7440-66-6	ZINC	ug/l	10.2 UJ	11.8 UJ	5910 J	9.49 UJ	30.8 UJ
57-12-5	CYANIDE	ug/l	3 U	123	224	27	5 U

			Dup of MW-8			
Con Ed - Lud		Location ID:	MW-18	MW-9	FIELDQC	FIELDQC
Validated Gro	undwater Analytical Data	Sample ID:	MW-18	MW-9	FB112712	TRIPBLANK
		Lab Sample Id:	D4947-03	D4947-04	D4947-11	D4947-10
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4947	D4947	D4947	D4947
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	11/26/2012 13:20	11/26/2012 14:45	11/27/2012 13:45	11/20/2012 9:00
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	VOLATILES		0.4.77	0.4.77	0.4.77	0.4.77
71-55-6	1,1,1-TRICHLOROETHANE	ug/l	0.4 U	0.4 U	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U
76-13-1 79-00-5	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
	1,1,2-TRICHLOROETHANE	ug/l	0.38 U	0.38 U	0.38 U 0.36 U	0.38 U 0.36 U
75-34-3 75-35-4	1,1-DICHLOROETHANE 1,1-DICHLOROETHENE	ug/l	0.36 U 0.47 U	0.36 U 0.47 U	0.36 U 0.47 U	0.36 U 0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/l ug/l	0.47 U 0.65 U	0.47 U 0.65 U	0.47 U 0.65 U	0.47 U 0.65 U
120-82-1	1,2,4-TRICHLOROBENZENE		0.62 U	0.62 U	0.62 U	0.62 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/l ug/l	0.46 U	0.62 U 0.46 U	0.46 U	0.46 U
106-93-4	1,2-DIBROMO-3-CHLOROPROPANE 1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/l ug/l	0.46 U 0.41 U	0.46 U 0.41 U	0.46 U 0.41 U	0.46 U 0.41 U
95-50-1	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE) 1,2-DICHLOROBENZENE	ug/l ug/l	0.41 U 0.45 U	0.41 U 0.45 U	0.41 U 0.45 U	0.41 U 0.45 U
107-06-2	1,2-DICHLOROBENZENE 1,2-DICHLOROETHANE	ug/l ug/l	0.48 U	0.43 U 0.48 U	0.43 U 0.48 U	0.48 U
78-87-5	1,2-DICHLOROPROPANE	ug/l ug/l	0.46 U	0.46 U	0.46 U	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/l ug/l	0.43 U	0.40 U	0.43 U	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/l	0.43 U	0.43 U	0.32 U	0.43 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/l	50 UJ	50 UJ	50 UJ	50 UJ
591-78-6	2-HEXANONE	ug/l	1.9 U	1.9 U	1.9 U	1.9 U
67-64-1	ACETONE	ug/l	2.8 UJ	2.8 UJ	2.8 UJ	2.8 UJ
71-43-2	BENZENE	ug/l	0.32 U	0.32 U	0.32 U	0.32 U
74-97-5	BROMOCHLOROMETHANE	ug/l	2.2 U	2.2 U	2.2 U	2.2 U
75-27-4	BROMODICHLOROMETHANE	ug/l	0.36 UJ	0.36 UJ	0.36 UJ	0.36 UJ
75-25-2	BROMOFORM	ug/l	0.47 UJ	0.47 UJ	0.47 UJ	0.47 UJ
74-83-9	BROMOMETHANE	ug/l	0.62 UJ	0.62 UJ	0.62 UJ	0.62 UJ
75-15-0	CARBON DISULFIDE	ug/l	0.54 U	0.54 U	2 J	0.54 U
56-23-5	CARBON TETRACHLORIDE	ug/l	0.62 UJ	0.62 UJ	0.62 UJ	0.62 UJ
108-90-7	CHLOROBENZENE	ug/l	0.49 U	0.49 U	0.49 U	0.49 U
75-00-3	CHLOROETHANE	ug/l	0.66 U	0.66 U	0.66 U	0.66 U
67-66-3	CHLOROFORM	ug/l	0.34 U	0.34 U	0.34 U	0.34 U
74-87-3	CHLOROMETHANE	ug/l	0.54 UJ	0.54 UJ	0.54 UJ	0.54 UJ
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U
110-82-7	CYCLOHEXANE	ug/l	0.55 U	0.55 U	0.55 U	0.55 U
124-48-1	DIBROMOCHLOROMETHANE	ug/l	0.52 UJ	0.52 UJ	0.52 UJ	0.52 UJ
75-71-8	DICHLORODIFLUOROMETHANE	ug/l	0.55 U	0.55 U	0.55 U	0.55 U
100-41-4	ETHYLBENZENE	ug/l	0.53 U	0.53 U	0.53 U	0.53 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/l	0.95 U	0.95 U	0.95 U	0.95 U
79-20-9	METHYL ACETATE	ug/l	0.83 UJ	0.83 UJ	0.83 UJ	0.83 UJ
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/l	1.3 U	1.3 U	1.3 U	1.3 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/l	2.1 UJ	2.1 UJ	2.1 UJ	2.1 UJ
108-87-2	METHYLCYCLOHEXANE	ug/l	0.68 U	0.68 U	0.68 U	0.68 U
75-09-2	METHYLENE CHLORIDE	ug/l	0.41 U	0.41 U	4.3 J	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/l	0.43 U	0.43 U	0.43 U	0.43 U
100-42-5	STYRENE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/l	0.49 J	1.4 J	0.35 U	0.35 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/l	0.27 U	0.27 U	0.27 U	0.27 U
108-88-3	TOLUENE TRANS 12 DIGH OROSTHENE	ug/l	0.37 U	0.37 U	0.37 U	0.37 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/l	0.29 UJ	0.29 UJ	0.29 UJ	0.29 UJ
79-01-6 75-69-4	TRICHLOROETHYLENE (TCE)	ug/l	0.28 U	0.28 U	0.28 U	0.28 U
	TRICHLOROFLUOROMETHANE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U
75-01-4	VINYL CHLORIDE	ug/l	0.34 U	0.34 U	0.34 U	0.34 U

			Dup of MW-8			
Con Ed - Ludlo		Location ID:	MW-18	MW-9	FIELDQC	FIELDQC
Validated Grou	ındwater Analytical Data	Sample ID:	MW-18	MW-9	FB112712	TRIPBLANK
		Lab Sample Id:	D4947-03	D4947-04	D4947-11	D4947-10
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4947	D4947	D4947	D4947
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	11/26/2012 13:20	11/26/2012 14:45	11/27/2012 13:45	11/20/2012 9:00
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
95-94-3	SEMIVOLATILES 1,2,4,5-TETRACHLOROBENZENE	ug/l	0.21 U	0.21 U	0.21 U	
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/l	0.21 U	0.21 UJ	0.21 U	
95-95-4	2,4,5-TRICHLOROPHENOL	ug/l	0.42 U	0.42 U	0.41 U	
88-06-2	2,4,6-TRICHLOROPHENOL	ug/l	0.58 U	0.59 U	0.58 U	
120-83-2	2,4-DICHLOROPHENOL	ug/l	0.69 U	0.69 U	0.68 U	
105-67-9	2,4-DIMETHYLPHENOL	ug/l	0.74 U	0.75 U	0.73 U	
51-28-5	2,4-DINITROPHENOL	ug/l	2.2 UJ	2.2 UJ	2.2 UJ	
121-14-2	2,4-DINITROTOLUENE	ug/l	1.1 U	1.1 U	1.1 U	
606-20-2	2,6-DINITROTOLUENE	ug/l	0.33 U	0.34 U	0.33 U	
91-58-7	2-CHLORONAPHTHALENE	ug/l	0.17 U	0.17 U	0.16 U	
95-57-8	2-CHLOROPHENOL	ug/l	0.56 U	0.57 U	0.56 U	
91-57-6	2-METHYLNAPHTHALENE	ug/l	0.33 U	0.34 U	0.33 U	
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/l	0.25 U	0.25 U	0.25 U	
88-74-4	2-NITROANILINE	ug/l	0.51 U	0.52 U	0.51 U	
88-75-5	2-NITROPHENOL	ug/l	0.54 U	0.55 U	0.54 U	
91-94-1	3,3'-DICHLOROBENZIDINE	ug/l	2.1 U	2.1 U	2.1 U	
	3+4-Methylphenols	ug/l	0.4 U	0.4 U	0.39 U	
99-09-2	3-NITROANILINE	ug/l	1.1 U	1.1 U	1.1 U	
534-52-1 101-55-3	4,6-DINITRO-2-METHYLPHENOL 4-BROMOPHENYL PHENYL ETHER	ug/l	0.77 UJ 0.24 U	0.78 UJ 0.24 U	0.76 UJ 0.24 U	
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/l ug/l	0.24 U 0.42 U	0.24 U 0.42 U	0.24 U 0.41 U	
106-47-8	4-CHLOROANILINE	ug/l	2.1 U	3 U	2.9 U	
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/l	0.22 U	0.22 U	0.22 U	
100-01-6	4-NITROANILINE	ug/l	1.4 U	1.4 U	1.4 U	
100-02-7	4-NITROPHENOL	ug/l	2.1 U	2.1 U	2.1 U	
83-32-9	ACENAPHTHENE	ug/l	0.22 U	0.22 U	0.22 U	
208-96-8	ACENAPHTHYLENE	ug/l	0.73 U	0.74 U	0.72 U	
98-86-2	ACETOPHENONE	ug/l	0.15 U	0.15 U	0.14 U	
120-12-7	ANTHRACENE	ug/l	0.17 U	0.17 U	0.16 U	
1912-24-9	ATRAZINE	ug/l	0.42 U	0.42 U	0.41 U	
100-52-7	BENZALDEHYDE	ug/l	0.8 U	0.81 U	0.79 U	
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.17 U	0.17 U	0.16 U	
50-32-8	BENZO(A)PYRENE	ug/l	0.15 U	0.15 U	0.14 U	
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3 U	0.31 U	0.3 U	
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.3 U	0.31 U	0.3 U	
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.19 U	0.19 U	0.19 U	
85-68-7 92-52-4	BENZYL BUTYL PHTHALATE BIPHENYL (DIPHENYL)	ug/l	0.2 U	0.2 U	0.2 U	
92-32- 4 111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/l	0.16 U 0.57 U	0.16 U 0.58 U	0.15 U 0.57 U	
111-44-4	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/l ug/l	0.57 U	0.58 U	0.57 U	
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/l ug/l	0.37 U 0.18 U	0.38 U 0.18 U	0.37 U 0.18 U	
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	0.17 U	0.17 U	0.16 U	
105-60-2	CAPROLACTAM	ug/l	2.1 U	2.1 U	2.1 U	
86-74-8	CARBAZOLE	ug/l	0.23 U	0.23 U	0.23 U	
218-01-9	CHRYSENE	ug/l	0.19 U	0.19 U	0.19 U	
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.44 U	0.44 U	0.43 U	
132-64-9	DIBENZOFURAN	ug/l	0.25 U	0.25 U	0.25 U	
84-66-2	DIETHYL PHTHALATE	ug/l	0.4 U	0.4 U	0.39 U	
131-11-3	DIMETHYL PHTHALATE	ug/l	0.23 U	0.23 U	0.23 U	
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	2.1 U	2.1 U	2.1 U	
117-84-0	DI-N-OCTYLPHTHALATE	ug/l	0.53 U	0.54 U	0.53 U	
206-44-0	FLUORANTHENE	ug/l	0.42 U	0.42 U	0.41 U	
86-73-7	FLUORENE	ug/l	0.32 U	0.33 U	0.32 U	
118-74-1	HEXACHLOROBENZENE	ug/l	0.19 U	0.19 U	0.19 U	
87-68-3	HEXACHLOROBUTADIENE	ug/l	0.26 U	0.26 U	0.26 U	
77-47-4 67-72-1	HEXACHLOROCYCLOPENTADIENE HEXACHLOROETHANE	ug/l	0.25 U	0.25 U	0.25 U	
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/l ug/l	0.26 U 0.16 U	0.26 U 0.16 U	0.26 U 0.15 U	
78-59-1	ISOPHORONE	ug/l ug/l	0.16 U 0.31 U	0.16 U 0.32 U	0.13 U 0.31 U	
91-20-3	NAPHTHALENE	ug/l	0.31 U 0.13 U	0.32 U 0.13 U	0.31 U 0.12 U	
91-20-3 98-95-3	NITROBENZENE	ug/l ug/l	0.13 U 0.71 U	0.13 U 0.72 U	0.12 U 0.7 U	
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/l	0.71 U 0.21 U	0.72 U 0.21 U	0.7 U 0.21 U	
86-30-6	N-NITROSODIPHENYLAMINE	ug/l	0.63 U	0.63 U	0.62 U	
87-86-5	PENTACHLOROPHENOL	ug/l	1.8 U	1.8 U	1.8 U	
85-01-8	PHENANTHRENE	ug/l	0.27 U	0.27 U	0.27 U	
03-01-0			=			
108-95-2	PHENOL	ug/l	0.22 U	0.22 U	0.22 U	

			Dup of MW-8			
Con Ed - Luc	flow Site	Location ID:	MW-18	MW-9	FIELDQC	FIELDQC
Validated Gr	oundwater Analytical Data	Sample ID:	MW-18	MW-9	FB112712	TRIPBLANK
	•	Lab Sample Id:	D4947-03	D4947-04	D4947-11	D4947-10
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	D4947	D4947	D4947	D4947
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	11/26/2012 13:20	11/26/2012 14:45	11/27/2012 13:45	11/20/2012 9:00
		Validated:	1/23/2013	1/23/2013	1/23/2013	1/23/2013
CAS NO.	COMPOUND	UNITS:				
	INORGANICS					
7429-90-5	ALUMINUM	ug/l	9290 J	26.5 UJ	61.4 J	
7440-36-0	ANTIMONY	ug/l	4 UJ	4 UJ	4 UJ	
7440-38-2	ARSENIC	ug/l	15.3 J	13.7	2.1 U	
7440-39-3	BARIUM	ug/l	169 J	478 J	2.22 J	
7440-41-7	BERYLLIUM	ug/l	0.38 J	0.35 UJ	0.35 UJ	
7440-43-9	CADMIUM	ug/l	0.87 J	0.27 J	0.25 U	
7440-70-2	CALCIUM	ug/l	90000 J	198000 J	295 J	
7440-47-3	CHROMIUM, TOTAL	ug/l	37.9 UJ	11.8 UJ	25.7 J	
7440-48-4	COBALT	ug/l	12.8 J	2.9 U	2.9 U	
7440-50-8	COPPER	ug/l	59 J	1.11 J	1 UJ	
7439-89-6	IRON	ug/l	18700 J	17400 J	161 J	
7439-92-1	LEAD	ug/l	85.2 J	1.3 U	1.3 U	
7439-95-4	MAGNESIUM	ug/l	31000 J	88100 J	53.7 J	
7439-96-5	MANGANESE	ug/l	2030 J	4210 J	4.96 J	
7439-97-6	MERCURY	ug/l	0.092 U	0.092 U	0.092 U	
7440-02-0	NICKEL	ug/l	31 UJ	10 U	10.5	
7440-09-7	POTASSIUM	ug/l	10900 J	22200 J	763 J	
7782-49-2	SELENIUM	ug/l	2.4 U	2.82 J	2.4 U	
7440-22-4	SILVER	ug/l	0.75 UJ	0.75 UJ	0.75 UJ	
7440-23-5	SODIUM	ug/l	677000 J	18600 UJ	17600 J	
7440-28-0	THALLIUM	ug/l	1.2 UJ	1.2 UJ	1.2 UJ	
7440-62-2	VANADIUM	ug/l	20.6 J	3.05 UJ	3.05 UJ	
7440-66-6	ZINC	ug/l	90.7 J	3.25 UJ	14 J	
57-12-5	CYANIDE	ug/l	5 U	8 U	3 J	

DATA USABILITY SUMMARY REPORT

LUDLOW FORMER MGP SITE

SUPPLEMENTAL REMEDIAL INVESTIGATION

Prepared For:



CONSOLIDATED EDISON COMPANY OF NEW YORK, INC.

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PARSONS

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APRIL 2013

TABLE OF CONTENTS

	Page
SECTION 1 DATA USABILITY SUMMARY	1-1
1.1 LABORATORY DATA PACKAGES	1-1
1.2 SAMPLING AND CHAIN-OF-CUSTODY	1-1
1.3 LABORATORY ANALYTICAL METHODS	1-1 1-2
SECTION 2 DATA VALIDATION REPORT	2-1
2.1 SOIL GAS SAMPLES	

LIST OF ATTACHMENTS

ATTACHMENT A – VALIDATED LABORATORY DATA

SECTION 1

DATA USABILITY SUMMARY

Soil gas samples were collected from the Consolidated Edison Ludlow Former MGP site on October 24, 2012. Analytical results from these samples were validated and reviewed by Parsons for usability with respect to the following requirements:

- Work Plan,
- NYSDEC Analytical Services Protocol (ASP), and
- USEPA Region II Standard Operating Proce dures (SOPs) for organic and inorganic data review.

The analytical laboratory for this project was Eurofins – Air Toxics (Eurofins). This laboratory is certified to perform project analyses through the New York State Department of Health (NYSDOH) Environmental Laboratory Approval Program (ELAP).

1.1 LABORATORY DATA PACKAGES

The laboratory data package turnaround time, defined as the time from sample receipt by the laboratory to receipt of the analytical data packages by Parsons, was 24 days for the project samples.

The data packages received from Eurofins were paginated, complete, and overall were of good quality. Comments on specific quality control (QC) and other requirements are discussed in detail in the attached data validation reports which are summarized in Section 2.

1.2 SAMPLING AND CHAIN-OF-CUSTODY

The samples were collected, properly preser ved, shipped under a chain-of-custody (COC) record, and received at Eurofins within one day of sampling. All samples were received intact and in good condition at the laboratory.

1.3 LABORATORY ANALYTICAL METHODS

The soil gas sam ples that were collected f rom the site were analyzed f or volatile organic compounds (VOCs). Summaries of issues concerning these laboratory analyses are presented in Subsections 1.3.1. The data qualifications resulting from the data validation review and statements on the laboratory analytical precision, accuracy, representativeness, completeness, and comparability (PARCC) are discussed for each analytical method in Section 2. The laboratory data were reviewed and may be qualified with the following validation flags:

- "U" not detected at the value given,
- "UJ" estimated and not detected at the value given,
 - "J" estimated at the value given,

"N" - presumptive evidence at the value given, and

"R" - unusable value.

The validated laboratory data were tabulated and are presented in Attachment A.

1.3.1 Volatile Organic Analysis

Soil gas sam ples were analyzed for VOCs us ing the USEPA TO-15 analytical m ethod. Certain reported results for the VOC sam ples were qualified as estimated based upon laboratory control sample (LCS) recoveries. The reporte d VOC analytical results were 100% com plete (i.e., usable) for the soil gas data. PARCC requirements were met.

SECTION 2

DATA VALIDATION REPORT

2.1 SOIL GAS SAMPLES

Data review has been completed for data packages generated by Eurofins containing analytical results from soil gas samples collected from the site. All of these samples were properly preserved, shipped under a COC record, and received intact by the analytical laboratory. Analytical results for these samples were contained with SDG 1210554. Data validation was performed for all samples in accordance with the most current editions of the USEPA Region II SOPs for organic and inorganic data review. This data validation and usability report is presented by analysis type. The validated laboratory data are presented in Attachment A.

2.1.1 Volatiles

The following items were reviewed for compliancy in the volatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- Matrix spike/matrix spike duplicate (MS/MSD) precision and accuracy
- Laboratory control sample (LCS) recoveries
- Laboratory method blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered com pliant and acceptable in accordance with the validation protocols with the exception of LCS recoveries as discussed below.

LCS Recoveries

All LCS recoveries were considered acceptable and within QC limits with the exception of the low LCS recoveries for ethanol (60%R, 61%R; QC limit 70-130%R) associated with all samples. Therefore, the ethanol results fo r these sam ples which were nondetects, were considered estimated, possibly biased low, and qualified "UJ" for the affected samples.

Usability

All soil gas sample results were considered usable following data validation.

Summary

The quality assurance objectives f or measurement data included considerations f or precision, accuracy, representativeness, completeness, and comparability. The volatile soil gas data presented by Eurofins were 100% complete (i.e., usable). The validated volatile laboratory data are tabulated and presented in Attachment A.

ATTACHMENT A VALIDATED LABORATORY DATA

C 1:1 + 1	E I.	I (ID	CVI 1	CV 2	CVIA
Consolidated		Location ID:	SV-1	SV-3	SV-4
Ludlow Stree	t Site Analytical Data	Sample ID: Lab Sample Id:	SV-1 1210554-03A	SV-3 1210554-01A	SV-4 1210554-02A
validated Air	Analytical Data				
		Source:	AIRTOXICS	AIRTOXICS	AIRTOXICS
		SDG:	1210554	1210554	1210554
		Matrix:	SOIL VAPOR	SOIL VAPOR	SOIL VAPOR
		Sampled:	10/24/2012	10/24/2012	10/24/2012
		Validated:	4/15/2013	4/15/2013	4/15/2013
CAS NO.	COMPOUND	UNITS:			
67.64.1	VOLATILES	_ , _	2.00	150	220
67-64-1	ACETONE	ug/m3	260	150	220
107-05-1	ALLYL CHLORIDE (3-CHLOROPROPENE)	ug/m3	11 U	11 U	11 U
100-44-7	ALPHA-CHLOROTOLUENE	ug/m3	4.6 U	4.4 U	4.5 U
71-43-2	BENZENE	ug/m3	20	2.7 U	8.6
75-27-4	BROMODICHLOROMETHANE	ug/m3	6 U	5.7 U	5.9 U
75-25-2	BROMOFORM	ug/m3	9.2 U	8.8 U	9 U
74-83-9	BROMOMETHANE	ug/m3	35 U	33 U	34 U
106-99-0 75-15-0	1,3-BUTADIENE	ug/m3	34 17	1.9 U	4.4 14
56-23-5	CARBON DISULFIDE	ug/m3	5.6 U	11 U	5.5 U
108-90-7	CARBON TETRACHLORIDE	ug/m3		5.4 U 3.9 U	3.3 U 4 U
	CHLOROBENZENE	ug/m3 ug/m3	4.1 U		
75-00-3 67-66-3	CHLOROETHANE CHLOROFORM	ug/m3 ug/m3	9.4 U 4.4 U	9 U 4.2 U	9.2 U 4.3 U
74-87-3	CHLOROMETHANE	ug/m3 ug/m3	4.4 U 18 U	4.2 U 18 U	4.3 U 18 U
110-82-7	CYCLOHEXANE	ug/m3	7.9	2.9 U	3 U
124-48-1	DIBROMOCHLOROMETHANE	ug/m3	7.9 7.6 U	7.3 U	7.4 U
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/m3	6.9 U	6.6 U	6.7 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/m3	5.9	4.2 U	4.3 U
75-34-3	1,1-DICHLOROETHANE	ug/m3	3.6 U	3.5 U	3.5 U
107-06-2	1,2-DICHLOROETHANE	ug/m3	3.6 U	3.5 U	3.5 U
75-35-4	1,1-DICHLOROETHENE	ug/m3	3.5 U	3.4 U	3.5 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/m3	3.5 U	3.4 U	3.5 U
156-60-5	TRANS-1,2-DICHLOROETH TLENE	ug/m3	3.5 U	3.4 U	3.5 U
95-50-1	1,2-DICHLOROBENZENE	ug/m3	5.4 U	5.1 U	5.3 U
541-73-1	1,3-DICHLOROBENZENE	ug/m3	5.4 U	5.1 U	5.3 U
106-46-7	1,4-DICHLOROBENZENE	ug/m3	5.4 U	5.1 U	5.3 U
78-87-5	1,2-DICHLOROPROPANE	ug/m3	4.1 U	4 U	4 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/m3	4.1 U	3.9 U	4 U
10061-01-3	TRANS-1,3-DICHLOROPROPENE	ug/m3	4.1 U	3.9 U	4 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/m3	13 U	12 U	13 U
64-17-5	ETHANOL	ug/m3	6.7 UJ	6.4 UJ	6.6 UJ
622-96-8	4-ETHYLTOLUENE	ug/m3	4.4 U	4.2 U	4.3 U
100-41-4	ETHYLBENZENE	ug/m3	5.4	3.7 U	5.8
76-14-2	FREON 114	ug/m3	6.2 U	6 U	6.1 U
142-82-5	HEPTANE	ug/m3	13	3.5 U	190
87-68-3	HEXACHLOROBUTADIENE	ug/m3	38 U	36 U	37 U
110-54-3	N-HEXANE	ug/m3	27	3 U	360
591-78-6	2-HEXANONE	ug/m3	15 U	14 U	14 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/m3	4.4 U	4.2 U	4.3 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/m3	11	10 U	10 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/m3	3.7 U	3.5 U	3.6 U
75-09-2	METHYLENE CHLORIDE	ug/m3	31 U	30 U	30 U
67-63-0	2-PROPANOL	ug/m3	8.8 U	8.4 U	8.6 U
103-65-1	N-PROPYLBENZENE	ug/m3	4.4 U	4.2 U	4.3 U
100-42-5	STYRENE	ug/m3	3.8 U	3.6 U	3.7 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/m3	3.2 U	3.1 U	3.2 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/m3	6.1 U	5.9 U	6 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/m3	28	100	24
109-99-9	TETRAHYDROFURAN	ug/m3	2.6 U	2.5 U	2.6 U
108-88-3	TOLUENE	ug/m3	35	23	60
120-82-1	1,2,4-TRICHLOROBENZENE	ug/m3	26 U	25 U	26 U
71-55-6	1,1,1-TRICHLOROETHANE	ug/m3	6.5	4.7 U	4.8 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/m3	4.9 U	4.7 U	4.8 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/m3	4.8 U	4.6 U	4.7 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/m3	6.8 U	6.6 U	6.7 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/m3	5 U	4.8 U	4.9 U
95-63-6	1,2,4-TRIMETHYLBENZENE	ug/m3	4.4 U	4.2 U	4.3 U
108-67-8	1,3,5-TRIMETHYLBENZENE (MESITYLENE)	ug/m3	4.4 U	4.2 U	4.3 U
540-84-1	2,2,4-TRIMETHYLPENTANE	ug/m3	12	4 U	4.1 U
75-01-4	VINYL CHLORIDE	ug/m3	2.3 U	2.2 U	2.2 U
179601-23-1	M,P-XYLENE	ug/m3	19	11	20
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/m3	5.2	3.7 U	7

DATA USABILITY SUMMARY REPORT

LUDLOW FORMER MGP SITE

SUPPLEMENTAL REMEDIAL INVESTIGATION

Prepared For:



CONSOLIDATED EDISON COMPANY OF NEW YORK, INC.

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TABLE OF CONTENTS

	<u>Page</u>
SECTION 1 DATA USABILITY SUMMARY	1-1
1.1 LABORATORY DATA PACKAGES	1-1
1.2 SAMPLING AND CHAIN-OF-CUSTODY	1-1
1.3 LABORATORY ANALYTICAL METHODS	1-1
1.3.1 Volatile Organic Analysis	1-2
1.3.2 Semivolatile Organic Analysis	
1.3.3 Inorganics Analysis	1-2
SECTION 2 DATA VALIDATION REPORT	2-1
2.1 GROUNDWATER	
2.1.1 Volatiles	2-1
2.1.2 Semivolatiles	2-3
2.1.3 Inorganics	2-5

LIST OT ATTACHMENTS

ATTACHMENT A VALIDATED LABORATORY DATA

SECTION 1

DATA USABILITY SUMMARY

Groundwater samples were collected from the Consolidated Edison Ludlow Street Site from April 4, 2013 through April 8, 2013. Analytical results from these samples were validated and reviewed by Parsons for usability with respect to the following requirements:

- Work Plan,
- NYSDEC Analytical Services Protocol (ASP), and
- USEPA Region II Standard Operating Proce dures (SOPs) for organic and inorganic data review.

The analytical laboratory for this project was Chemtech. This laboratory is certified to perform project analyses through the New York State Department of Health (NYSDOH) Environmental Laboratory Approval Program (ELAP).

1.1 LABORATORY DATA PACKAGES

The laboratory data package turnaround time, defined as the time from sample receipt by the laboratory to receipt of the analytical data packages by Parsons, was 14 days for the project samples.

The data packages received from Chemtech were paginated, complete, and overall were of good quality. Comments on specific quality control (QC) and other requirements are discussed in detail in the attached data validation report which is summarized by media in Section 2.

1.2 SAMPLING AND CHAIN-OF-CUSTODY

The samples were collected, properly preser ved, shipped under a chain-of-custody (COC) record, and received at Chem tech within one to four days of sam pling. All sam ples were received intact and in good condition at Chemtech.

1.3 LABORATORY ANALYTICAL METHODS

The groundwater sam ples that were collected from the site were analyzed for volatile organic compounds (VOCs), semivolatile organic compounds (SVOCs), metals, and cyanide. Summaries of issues concerning these laborator y analyses are presented in Subsections 1.3.1 through 1.3.3. The data qualifications resulting from the data validation review and statem ents on the laboratory analytical precision, accurac y, representativeness, com pleteness, and comparability (PARCC) are discussed for each analytical m ethod in Section 2. The laboratory data were reviewed and may be qualified with the following validation flags:

"U" - not detected at the value given,

"UJ" - estimated and not detected at the value given,

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"J" - estimated at the value given,

"N" - presumptive evidence at the value given, and

"R" - unusable value.

The validated laboratory data were tabulated and are presented in Attachment A.

1.3.1 Volatile Organic Analysis

Groundwater samples were analyzed for VOCs using the USEPA SW-846 8260C analytical method. Certain reported results for the VOC samples were qualified as estim ated based upon matrix spike/matrix spike duplicate (MS/MSD) recoveries and instrument calibrations. The reported VOC analytical results were 100% complete (i.e., usable) for the groundwater data. PARCC requirements were met.

1.3.2 Semivolatile Organic Analysis

Groundwater samples were analyzed for SVOCs using the USEPA SW -846 8270D analytical method. Certain reported results for the SVOC samples were qualified as estim ated based upon MS/MSD recoveries, instrum ent calibrations, internal standard responses, and field duplicate precision. Certain reported results for the SVOC samples were considered unusable and qualified "R" based upon poor MS/MSD recoveri es. The reported SVOC analytical results were 99.7% complete (i.e., usable) for the gr oundwater data. PARCC requirem ents were met overall.

1.3.3 Inorganics Analysis

Groundwater samples were analyzed for metals and cyanide using the USEPA SW -846 6010B/7470A/7471A/9012B analytical methods. Certain reported results for the inorganics samples were qualified as estimated based upon matrix spike recoveries, serial dilutions, and field duplicate precision. Certain reported results for the inorganic samples were considered unusable and qualified "R" based upon poor matrix spike recoveries. The reported inorganic analytical results were considered 97.2% complete (i.e., usable) for the groundwater data. PARCC requirements were met overall.

SECTION 2

DATA VALIDATION REPORT

2.1 GROUNDWATER

Data review has been completed for data packages generated by Chem tech containing groundwater samples collected from the site. All of these samples were properly preserved, shipped under a COC record, and received intact by the analytical laboratory. The analytical results were presented by the laboratory in one sample delivery group (SDG): E1768. Data validation was performed for all samples in accordance with the most current editions of the USEPA Region II SOPs for organic and inorganic containing the validated laboratory data are presented in Attachment A.

2.1.1 Volatiles

The following items were reviewed for compliancy in the volatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- Matrix spike/matrix spike duplicate (MS/MSD) precision and accuracy
- Laboratory control sample (LCS) recoveries
- Laboratory method blank and field equipment/trip blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, blank contamination, and initial and continuing calibrations as discussed below.

MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent dif ference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC lim its for designated project spiked samples with the exception of the high MS/MSD accuracy results for chlorom ethane,

methyl tert-butyl ether, 1,1-dichloroethane, 2-butanone, bromochloromethane, chloroform, 1,1,1-trichloroethane, 4-methyl-2-pentanone, 2-hexanone, 1,2-dibromoethane, bromoform, 1,1,2,2-tetrachloroethane, 1,2-dibromo-3-chloropropane, and 1,4-dioxane; and the low MS/MSD accuracy results for bromomethane and methylcyclohexane during the spiked analyses of sample MW-5. The positive 2-butanone and 4-m ethyl-2-pentanone results for the parent sam ple were considered estimated, possibly biased hi gh, and qualified "J" based upon high MS/MSD accuracy results. The nondetected brom omethane and methylcyclohexane results for the parent sample were considered estim ated, possibly biased low, and qualified "UJ" based upon low MS/MSD accuracy results. The remaining noncompliant compounds did not required validation qualification for the parent sample.

LCS Recoveries

All LCS recoveries were considered accepta ble and within QC lim its for all com pounds with the exception of the high LCS recovery for 1,2,3-trichlorobenzene (136%R; 58-130%R) associated with sam ple MW-5DL. Validati on qualification was not required since this compound was not detected in this sample.

Blank Contamination

The field equipm ent blank FB040813 associat ed with groundwater sam ples contained carbon disulfide at a concentration of $5.4 \mu g/L$. Since this compound was not detected in the project samples, validation qualification was not required.

Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative st andard deviation (%RSD) of 20% with the exception of acetone (20.3%RSD) and chloroet hane (20.3%RSD) in the initial calibration associated with samples MW-5 and -5DL; bromomethane (22.1%RSD), acetone (23.7%RSD), and cyclohexane (42.6%RSD) in the initial calibration associated with all samples except MW-5, -5DL, and -3ADL; and dichlorodifluorom ethane (25.4%RSD), chlorom ethane (24.8%RSD), chloroethane (23.3%RSD), 1,1,2-trichlorotrifluor oethane (20.2%RSD), cis-1,3-dichloropropene (21.6%RSD), 2-hexanone (27.9%RSD), dibrom ochloromethane (27.3%RSD), brom oform (32.5%RSD), and 1,2-dibromo-3-chloropropane (23.6%RSD) in the initial calibration associated with sample MW-3ADL. The results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within + _20% with the exception of m _ ethyl acetate (-21.25%D) and acetone (-24.14%D) in the continuing _ calibration associated with sam ple MW-5DL; cyclohexane (-20.59%D) in the continuing ca libration associated with all sam ples except MW-5, _-5DL, and _-3ADL; and carbon tetrach _ loride (24.61%D), brom odichloromethane (28.04%D), _4-methyl-2-pentanone (29.2%D), tr ans-1,3-dichloropropene (35.28%D), cis-1,3-dichloropropene (26.97%D), 2-hexanone (35.74%D), dibromochloromethane (35.84%D), styrene (22.09%D), brom oform (34.04%D), and 1,4-dioxane (25%D) in the continuing calibration associated with MW-3ADL. Therefore, the sample results for these compounds were

considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

<u>Usability</u>

All volatile groundwater sample results were considered usable following data validation.

Summary

The quality assurance objectives f or measurement data included considerations f or precision, accuracy, representativeness, com pleteness, and com parability. The volatile groundwater data presented by Chem tech were 100% complete (i.e., usable). The validated volatile laboratory data are tabulated and presented in Attachment A.

2.1.2 Semivolatiles

The following items were reviewed for compliancy in the semivolatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- MS/MSD precision and accuracy
- LCS recoveries
- Laboratory method blank and field equipment blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered com pliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, initial and continuing calibrations, internal standard res ponses, and field duplicate precision as discussed below.

MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated spiked project samples with the exception of the low MS/MSD accuracy results for naphthalene and 4-nitroaniline; the less than 10% MS/MSD accuracy results for 3-nitroaniline and 3,3-dichlorobenzidine; and the high MS/MSD accuracy result for indeno(1,2,3-cd)pyrene during the

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spiked analyses of sample MW-5. Therefore, results for those compounds that fell below the QC limit were considered estim ated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the parent sample. However, the nondetected 3-nitroaniline and 3,3-dichlorobenzidine results we re considered unusable and qualified "R" for the parent sample. Validation qualification of indeno(1,2,3-cd)pyrene was not required since this compound was not detected in the parent sample.

LCS Recoveries

All LCS recoveries were considered acceptable and within QC limits with the exception of the high LCS recovery for indeno(1,2,3-cd)pyr ene (139%R; QC lim it 45-125%R) associated with all groundwater sam ples. Validation qualification of the groundwater sam ples was not required since this compound was not detected.

Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a m aximum percent relative st andard deviation (%RSD) of 20% with the exception of hexachlorocyclopentadiene (24.2%RSD), 2.4-dinitrophenol (47.3%RSD), 4nitrophenol (20.3%RSD), pentachlorophenol (20.7%RSD), and 4,6-dinitro-2-m ethylphenol (32.1%RSD) in the initial calibration associ ated with sam ples MW-1, -5, and -9; 2,4dinitrophenol (41.9%RSD), 4,6-dinitro-2-m ethylphenol (24.2%RSD), and pentachlorophenol (23.7%RSD) in the initial calibration associated with samples FB040813, MW-2, -3, -3A, -5DL, -7, and -4; hexachlorocyclopentadiene (36.7%RSD), 2,4-dinitrophenol (47.6%RSD), 4nitrophenol (32.9%RSD), and 4.6-dinitro-2-m ethylphenol (28%RSD) in the initial calibration associated with sam ples MW-7DL, -7DL2, -3ADL, and -3ADL2; and benzaldehyde (22.5%RSD), hexachlorocyclopentadiene (24.9%RSD), 2,4-dinitrophenol (41.1%RSD), and 4nitrophenol (21.0%RSD). The results for these compounds were considered estim ated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within + 20% with the exception of indeno(1,2,3-cd)pyrene (78.9%D), dibenzo(a,h)anthracene (36.9%D), and benzo(g,h,i)perylene (35.7%D) in the continuing calibration associated with sam ples MW-1, -5, and -9; 2,4-dinitrophenol (20.6%D) and acetophenone (22.8%D) in the continuing calibration associated with samples MW-2, -3, -3A, -5DL, and -7; 2,4-dinitrophenol (-25.4%D) in the continuing calibration associated with sample MW-4; and 4-nitrophenol (-30.2% D), di-n-octylphthalate (-23.5%D), dibenzo(a,h)anthracene (43.1%D), and benzo(g,h,i)perylene (49.2%D) in the continuing calibration of MW-3DL and -3ADL3. Therefore, the sam ple results for these com pounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

Internal Standard Responses

All internal standard (IS) responses and rete $\,$ ntion times were within specified QC ranges based on associated calibration standards (i.e., sa $\,$ mple's area count within -50% to +100% and retention times within ± 0.5 minutes of the standard) with the exception of the low responses for the IS chrysene-d12 in samples MW-7DL2, -3ADL2, -7DL, -3ADL, -3DL, and -3ADL3; and the

low responses for the IS perylene-d12 in sam ples MW-3DL and -3ADL3. Validation qualification of these samples was not required since final data results were not affected.

Field Duplicate Precision

All field duplicate precision results were considered acceptable with the exception of the precision for acenaphthylene (53%RPD) associated with sample MW-3 and its field duplicate MW-3A. Therefore, the acenaphthylene results for these samples were considered estimated and qualified "J".

Usability

All semivolatile groundwater sam ple results were considered usable following data validation with the exception of certain nondetected results based upon poor MS/MSD recoveries.

Summary

The quality assurance objectives f or measurement data included considerations f or precision, accuracy, representativeness, com pleteness, and com parability. The sem ivolatile groundwater data presented by Chem tech were 99.7% complete (i.e., usable). The validated semivolatile laboratory data are tabulated and presented in Attachment A.

2.1.3 Inorganics

The following items were reviewed for compliancy in the inorganics analysis:

- Custody documentation
- Holding times
- Initial and continuing calibration verifications
- Initial and continuing calibration blank, and laboratory preparation blank, and field equipment blank contamination
- Inductively coupled plasma (ICP) interference check sample (ICS)
- Matrix spike (MS) recoveries
- Laboratory duplicate precision
- Field duplicate precision
- Laboratory control sample (LCS) recoveries
- ICP serial dilutions
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered com pliant and acceptable in accordance with the validation protocols with the exception of blank contam ination, matrix spike recoveries, field duplicate precision, and serial dilutions as discussed below.

Blank Contamination

The field equipment blank FB040813 associated with the groundwater sam ples contained dissolved sodium and dissolved zinc at concentrations of 3160 and 26.9 μ g/L, respectively. Therefore, all results for these analytes less than the validation action concentrations were considered not detected and qualified "U" for the affected samples.

Matrix Spike Recoveries

All matrix spike recoveries for all analytes were considered acceptable and within the 75-125%R QC limit with the exception of the matrix spike recoveries for aluminum (165.8%R, 146.7%R), cadmium (18.9%R, 17.7%R), chromium (50.1%R, 47.5%R), cobalt (72.3%R, 68.6%R), silver (49%R, 42.5%R), thallium (61.5%R, 57.8%R), copper (52.6%R, 49.6%R), iron (133%R), lead (51.9%R, 45.2%R), mercury (71.1%R, 72%R), nickel (63.9%R, 61%R), vanadium (62%R, 58.4%R), and zinc (55.6%R, 47.6%R) associated with all samples. Therefore, positive results for those analytes where matrix spike recoveries exceeded the QC limit were considered estimated, possibly biased high, and qualified "J" for the affected samples. Results for those analytes where matrix spike recoveries fell below the QC limit were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples. However, nondetected cadmium results were considered unusable and qualified "R" for the affected samples based upon less 30% matrix spike recoveries.

Field Duplicate Precision

All field duplicate precision results were considered acceptable for the field duplicate pair MW-3 and MW-3A with the exception of the precision for cyanide (129%RPD). Therefore, the cyanide results were considered estimated and qualified "J" for the affected parent sample and its field duplicate.

ICP Serial Dilutions

All serial dilution results for designated project samples were considered acceptable with a percent difference (%D) less than 10% for all IC P analytes with the exception of calcium (13.7%D), lead (11.7%D), magnesium (18.7%D), manganese (12.9%D), and sodium (16.3%D) associated with all sam ples. Therefore, positive results for these analytes were considered estimated and qualified "J" for the affected samples.

Usability

All inorganics groundwater sample results were considered usable following data validation with the exception of certain nondetected results based upon poor matrix spike recoveries.

Summary

The quality assurance objectives f or measurement data included considerations f or precision, accuracy, representativeness, com pleteness, and com parability. The inorganics groundwater data presented by Chem tech were 97.2% complete (i.e., usable). The validated groundwater inorganics laboratory data are tabulated and presented in Attachment A.

ATTACHMENT A VALIDATED LABORATORY DATA

						Dup of MW-3
Consolidated I	Edison	Location ID:	MW-1	MW-2	MW-3	MW-3A
Ludlow Street		Sample ID:	MW-1-20130405	MW-2-20130408	MW-3-20130408	MW-3A-20130408
	undwater Analytical Data	Lab Sample Id:	E1768-13	E1768-07	E1768-08	E1768-09
SDG: E1768	andwater rinarytical Data	Source:	CTECH	CTECH	CTECH	CTECH
BDG. E1700		SDG:	E1768	E1768	E1768	E1768
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	4/5/2013 13:00	4/8/2013 8:22	4/8/2013 9:54	4/8/2013 9:54
		Validated:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
CAS NO.	VOLATILES	CIVIIS.				
71-55-6	1,1,1-TRICHLOROETHANE	ug/l	0.4 U	0.4 U	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/l	0.4 U	0.4 U	0.4 U	0.4 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/l	0.43 U	0.43 U	0.38 U	0.43 U
75-34-3	1,1-DICHLOROETHANE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U
75-35-4	1,1-DICHLOROETHENE	ug/l	0.47 U	0.47 U	0.47 U	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/l	0.47 U	0.47 U	0.47 U	0.47 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	-	0.46 U	0.46 U	0.46 U	0.46 U
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/l ug/l	0.46 U 0.41 U	0.46 U 0.41 U	0.46 U 0.41 U	0.46 U 0.41 U
95-50-1	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE) 1,2-DICHLOROBENZENE	ug/l ug/l	0.41 U 0.45 U	0.41 U 0.45 U	0.41 U 0.45 U	0.41 U 0.45 U
107-06-2	1,2-DICHLOROETHANE		0.48 U	0.48 U	0.48 U	0.43 U 0.48 U
78-87-5	1,2-DICHLOROPROPANE	ug/l	0.46 U	0.46 U	0.46 U	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/l	0.40 U	0.40 U	0.43 U	0.40 U
106-46-7	1,4-DICHLOROBENZENE	ug/l	0.43 U	0.43 U	0.43 U	0.43 U
591-78-6	2-HEXANONE	ug/l ug/l	1.9 U	1.9 U	1.9 U	1.9 U
67-64-1	ACETONE		7.8 J	8.7 J	0.5 UJ	0.5 UJ
71-43-2	BENZENE	ug/l	0.32 U	0.32 U	360	460
74-97-5	BROMOCHLOROMETHANE	ug/l	0.32 U	0.32 U	0.2 U	0.2 U
75-27-4	BROMODICHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U 0.36 U
75-25-2	BROMOFORM	ug/l	0.47 U	0.47 U	0.47 U	0.47 U
74-83-9	BROMOMETHANE	ug/l ug/l	0.47 U 0.2 UJ	0.47 U 0.2 UJ	0.47 U 0.2 UJ	0.47 U 0.2 UJ
75-15-0	CARBON DISULFIDE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
108-90-7	CHLOROBENZENE	ug/l	0.49 U	0.49 U	0.49 U	0.49 U
75-00-3	CHLOROETHANE	ug/l	0.2 U	0.2 U	0.49 U	0.2 U
67-66-3	CHLOROFORM	ug/l	0.34 U	0.34 U	0.34 U	0.34 U
74-87-3	CHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/l	0.35 U	0.49 J	0.35 U	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U
110-82-7	CYCLOHEXANE	ug/l	0.2 UJ	0.2 UJ	1.5 J	2.1 J
124-48-1	DIBROMOCHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
100-41-4	ETHYLBENZENE	ug/l	0.2 U	0.2 U	160	180
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/l	0.45 U	0.45 U	3.9 J	4.6 J
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/l	0.95 U	0.95 U	300	370
79-20-9	METHYL ACETATE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/l	1.3 U	1.3 U	1.3 U	1.3 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/l	2.1 U	2.1 U	2.1 U	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/l	0.2 U	0.2 U	3.1 J	4.2 J
75-09-2	METHYLENE CHLORIDE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/l	0.43 U	0.43 U	160	180
100-42-5	STYRENE	ug/l	0.36 U	0.36 U	41.5	42.6
1634-04-4	TERT-BUTYL METHYL ETHER	ug/l	0.35 U	0.35 U	1.7 J	1.8 J
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/l	0.27 U	0.27 U	0.27 U	0.27 U
108-88-3	TOLUENE	ug/l	0.37 U	0.37 U	260	310
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/l	0.29 U	0.29 U	0.29 U	0.29 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/l	0.28 U	0.28 U	0.28 U	0.28 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U
75-01-4	VINYL CHLORIDE	ug/l	0.34 U	0.34 U	0.34 U	0.34 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/l	100 U	100 U	100 U	100 U

						Dup of MW-3
Consolidated E	dison	Location ID:	MW-1	MW-2	MW-3	MW-3A
Ludlow Street S		Sample ID:	MW-1-20130405	MW-2-20130408	MW-3-20130408	MW-3A-20130408
	ındwater Analytical Data	Lab Sample Id:	E1768-13	E1768-07	E1768-08	E1768-09
SDG: E1768		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	E1768	E1768	E1768	E1768
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	4/5/2013 13:00	4/8/2013 8:22	4/8/2013 9:54	4/8/2013 9:54
		Validated:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:				
95-94-3	SEMIVOLATILES		0.22 U	0.202 U	0.204.11	0.202.11
58-90-2	1,2,4,5-TETRACHLOROBENZENE 2,3,4,6-TETRACHLOROPHENOL	ug/l	0.22 U 0.22 U	0.202 U 0.202 U	0.204 U 0.204 U	0.202 U 0.202 U
95-95-4	2.4.5-TRICHLOROPHENOL	ug/l ug/l	0.45 U	0.404 U	0.408 U	0.404 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/l	0.43 U	0.566 U	0.571 U	0.566 U
120-83-2	2,4-DICHLOROPHENOL	ug/l	0.74 U	0.667 U	0.673 U	0.667 U
105-67-9	2,4-DIMETHYLPHENOL	ug/l	0.8 U	0.717 U	0.724 U	0.717 U
51-28-5	2,4-DINITROPHENOL	ug/l	2.4 UJ	2.1 UJ	2.1 UJ	2.1 UJ
121-14-2	2,4-DINITROTOLUENE	ug/l	1.2 U	1 U	1.1 U	1 U
606-20-2	2,6-DINITROTOLUENE	ug/l	0.36 U	0.323 U	0.327 U	0.323 U
91-58-7	2-CHLORONAPHTHALENE	ug/l	0.18 U	0.162 U	0.163 U	0.162 U
95-57-8	2-CHLOROPHENOL	ug/l	0.61 U	0.545 U	0.551 U	0.545 U
91-57-6	2-METHYLNAPHTHALENE	ug/l	0.36 U	0.323 U	630	630
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/l	0.27 U	0.242 U	0.245 U	0.242 U
88-74-4	2-NITROANILINE	ug/l	0.55 U	0.495 U	0.5 U	0.495 U
88-75-5	2-NITROPHENOL	ug/l	0.58 U	0.525 U	0.531 U	0.525 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/l	2.2 U	2 U	2 U	2 U
	3+4-Methylphenols	ug/l	0.43 U	0.384 U	0.388 U	0.384 U
99-09-2	3-NITROANILINE	ug/l	1.2 U	1.1 U	1.1 U	1.1 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/l	0.83 UJ	0.747 UJ	0.755 UJ	0.747 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/l	0.26 U	0.232 U	0.235 U	0.232 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/l	0.45 U	0.404 U	0.408 U	0.404 U
106-47-8	4-CHLOROANILINE	ug/l	3.2 U	2.9 U	2.9 U	2.9 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/l	0.24 U	0.212 U	0.214 U	0.212 U
100-01-6	4-NITROANILINE	ug/l	1.5 U	1.4 U	1.4 U	1.4 U
100-02-7	4-NITROPHENOL	ug/l	2.2 U	2 U	2 U	2 U
83-32-9 208-96-8	ACENAPHTHENE	ug/l	0.24 U 0.79 U	0.212 U	22.5	32.2 130 J
98-86-2	ACENAPHTHYLENE ACETOPHENONE	ug/l	0.79 U 0.16 U	0.707 U 0.141 UJ	75.6 J 4.8 J	8 J
120-12-7	ANTHRACENE	ug/l ug/l	0.18 U	0.162 U	21.9	27.4
1912-24-9	ATRAZINE	ug/l	0.45 U	0.404 U	0.408 U	0.404 U
100-52-7	BENZALDEHYDE	ug/l	0.43 U	0.778 U	0.786 U	0.778 U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.18 U	0.162 U	12.1	12.2
50-32-8	BENZO(A)PYRENE	ug/l	0.16 U	0.141 U	8.5 J	8.5 J
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.33 U	0.293 U	6.4 J	6.5 J
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.33 UJ	0.293 U	0.296 U	0.293 U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.2 U	0.182 U	0.184 U	0.182 U
85-68-7	BENZYL BUTYL PHTHALATE	ug/l	0.21 U	0.192 U	0.194 U	0.192 U
92-52-4	BIPHENYL (DIPHENYL)	ug/l	0.17 U	0.152 U	25.2	36.2
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/l	0.62 U	0.556 U	0.561 U	0.556 U
111-44-4	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/l	0.62 U	0.556 U	0.561 U	0.556 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/l	0.19 U	0.172 U	0.173 U	0.172 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	0.18 U	0.162 U	0.163 U	0.162 U
105-60-2	CAPROLACTAM	ug/l	2.2 U	2 U	2 U	2 U
86-74-8	CARBAZOLE	ug/l	0.25 U	0.222 U	5.3 J	8.4 J
218-01-9	CHRYSENE DIRECTOR HOANTHID A CENTE	ug/l	0.2 U	0.182 U	12.3	12.2
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.47 UJ	0.424 U	0.429 U	0.424 U
132-64-9	DIBENZOFURAN	ug/l	0.27 U	0.242 U	6.1 J	8.9 J 0.384 U
84-66-2 131-11-3	DIETHYL PHTHALATE	ug/l	0.43 U	0.384 U	0.388 U	
84-74-2	DIMETHYL PHTHALATE DI-N-BUTYL PHTHALATE	ug/l	0.25 U 2.2 U	0.222 U	0.224 U	0.222 U
84-74-2 117-84-0	DI-N-OCTYLPHTHALATE	ug/l ug/l	0.57 U	2 U 0.515 U	2 U 0.52 U	2 U 0.515 U
206-44-0	FLUORANTHENE	ug/I ug/I	0.45 U	0.404 U	17.8	19.9
86-73-7	FLUORENE	ug/l	0.45 U	0.404 U 0.313 U	47.3	65.1
118-74-1	HEXACHLOROBENZENE	ug/l	0.33 U 0.2 U	0.182 U	0.184 U	0.182 U
87-68-3	HEXACHLOROBUTADIENE	ug/l	0.28 U	0.253 U	0.255 U	0.253 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/l	0.27 UJ	0.242 U	0.245 U	0.242 U
67-72-1	HEXACHLOROETHANE	ug/l	0.28 U	0.253 U	0.255 U	0.253 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/l	0.17 UJ	0.152 U	0.153 U	0.152 U
78-59-1	ISOPHORONE	ug/l	0.34 U	0.303 U	0.306 U	0.303 U
91-20-3	NAPHTHALENE	ug/l	0.13 U	0.121 U	2300	3900
98-95-3	NITROBENZENE	ug/l	0.76 U	0.687 U	0.694 U	0.687 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/l	0.22 U	0.202 U	0.204 U	0.202 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/l	0.67 U	0.606 U	0.612 U	0.606 U
87-86-5	PENTACHLOROPHENOL	ug/l	1.9 UJ	1.7 UJ	1.8 UJ	1.7 UJ
85-01-8	PHENANTHRENE	ug/l	0.29 U	0.263 U	72.9	110
108-95-2	PHENOL	ug/l	0.24 U	0.212 U	0.214 U	0.212 U
129-00-0	PYRENE	ug/l	0.22 U	0.202 U	28.4	31.6

Page 2 of 9

						Dup of MW-3
Consolidated	Edison	Location ID:	MW-1	MW-2	MW-3	MW-3A
Ludlow Stree		Sample ID:	MW-1-20130405	MW-2-20130408	MW-3-20130408	MW-3A-20130408
	oundwater Analytical Data	Lab Sample Id:	E1768-13	E1768-07	E1768-08	E1768-09
SDG: E1768		Source:	CTECH	CTECH	CTECH	CTECH
3DG. E1/06						
		SDG:	E1768	E1768	E1768	E1768
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	4/5/2013 13:00	4/8/2013 8:22	4/8/2013 9:54	4/8/2013 9:54
	T	Validated:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:				
	TOTAL METALS					
7429-90-5	ALUMINUM	ug/l	2380 J	368 J	96.1 J	81.4 J
7440-36-0	ANTIMONY	ug/l	8 U	9.48 J	8 U	8 U
7440-38-2	ARSENIC	ug/l	7.8 J	4.2 U	4.2 U	4.25 J
7440-39-3	BARIUM	ug/l	129	1990	134	128
7440-41-7	BERYLLIUM	ug/l	0.7 U	0.7 U	0.7 U	0.7 U
7440-43-9	CADMIUM	ug/l	R	1.22 J	0.52 J	R
7440-70-2	CALCIUM	ug/l	59000 J	262000 J	81900 J	77000 J
7440-47-3	CHROMIUM, TOTAL	ug/l	37.6 J	1.1 UJ	1.1 UJ	1.1 UJ
7440-48-4	COBALT	ug/l	5.8 UJ	6.91 J	5.8 UJ	5.8 UJ
7440-50-8	COPPER	ug/l	7.59 J	27.4 J	7.59 J	6.54 J
7439-89-6	IRON	ug/l	8770	6860	1600	1590
7439-92-1	LEAD	ug/l	6.36 J	10.1 J	40.3 J	37.6 J
7439-95-4	MAGNESIUM	ug/l	13400 J	33200 J	22600 J	21000 J
7439-96-5	MANGANESE	ug/l	2200 J	2160 J	1500 J	1400 J
7439-90-3	MERCURY		0.092 U	0.092 U	0.092 U	0.092 U
7440-02-0	NICKEL	ug/l	20.9 J	4.2 UJ	5.62 J	5.4 J
7440-02-0		ug/l	6390			
	POTASSIUM	ug/l		113000	15500	14300
7782-49-2	SELENIUM	ug/l	4.8 U	4.8 U	4.8 U	4.8 U
7440-22-4	SILVER	ug/l	1.5 UJ	1.5 UJ	1.5 UJ	1.5 UJ
7440-23-5	SODIUM	ug/l	272000 J	39700000 J	4620000 J	5420000 J
7440-28-0	THALLIUM	ug/l	2.4 UJ	2.4 UJ	2.4 UJ	2.4 UJ
7440-62-2	VANADIUM	ug/l	6.1 UJ	6.1 UJ	6.1 UJ	6.1 UJ
7440-66-6	ZINC	ug/l	18.4 J	35.4 J	20.8 J	21.1 J
	DISSOLVED METALS	1				
7429-90-5	ALUMINUM	ug/l	6.5 UJ			
7440-36-0	ANTIMONY	ug/l	8 U			
7440-38-2	ARSENIC	ug/l	4.2 U			
7440-39-3	BARIUM	ug/l	64.9			
7440-41-7	BERYLLIUM	ug/l	0.7 U			
7440-43-9	CADMIUM	ug/l	R			
7440-70-2	CALCIUM	ug/l	54500 J			
7440-47-3	CHROMIUM, TOTAL	ug/l	1.1 UJ			
7440-48-4	COBALT	ug/l	5.8 UJ			
7440-50-8	COPPER	ug/l	2 UJ			
7439-89-6	IRON	ug/l	20.4 U			
7439-92-1	LEAD	ug/l	2.6 UJ			
7439-95-4	MAGNESIUM	ug/l	10700 J			
7439-96-5	MANGANESE	ug/l	680 J			
7439-97-6	MERCURY	ug/l	0.092 U			
7440-02-0	NICKEL	ug/l	4.2 UJ			
7440-02-0	POTASSIUM		6360			
7782-49-2	SELENIUM	ug/l				
		ug/l	4.8 U			
7440-22-4	SILVER	ug/l	1.5 UJ			
7440-23-5	SODIUM	ug/l	271000 J			
7440-28-0	THALLIUM	ug/l	2.4 UJ			
7440-62-2	VANADIUM	ug/l	6.1 UJ			
7440-66-6	ZINC	ug/l	9.09 J			
1	OTHER	<u> </u>				
57-12-5	CYANIDE	ug/l	3 U	224	36 J	166 J

Consolidated	Edison	Location ID:	MW-4	MW-5	MW-7	MW-9
Ludlow Street		Sample ID:	MW-4-20130408	MW-5-20130405	MW-7-20130404	MW-9-20130405
	oundwater Analytical Data	Lab Sample Id:	E1768-06	E1768-03	E1768-12	E1768-01
SDG: E1768	And water I mary tear Bata	Source:	CTECH	CTECH	CTECH	CTECH
BDG. E1700		SDG:	E1768	E1768	E1768	E1768
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	4/8/2013 8:40	4/5/2013 14:10	4/4/2013 12:32	4/5/2013 10:35
		Validated:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/l	0.4 U	0.4 U	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/l	0.38 U	0.38 U	0.38 U	0.38 U
75-34-3	1,1-DICHLOROETHANE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U
75-35-4	1,1-DICHLOROETHENE	ug/l	0.47 U	0.47 U	0.47 U	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/l	0.46 U	0.46 U	0.46 U	0.46 U
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
95-50-1	1,2-DICHLOROBENZENE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
107-06-2	1,2-DICHLOROETHANE	ug/l	0.48 U	0.48 U	0.48 U	0.48 U
78-87-5	1,2-DICHLOROPROPANE	ug/l	0.46 U	0.46 U	0.46 U	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/l	0.43 U	0.43 U	0.43 U	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/l	0.32 U	0.32 U	0.32 U	0.32 U
591-78-6	2-HEXANONE	ug/l	1.9 U	1.9 U	1.9 U	1.9 U
67-64-1	ACETONE	ug/l	81.1 J	320 J	0.5 UJ	0.5 UJ
71-43-2	BENZENE	ug/l	0.32 U	35.8	1200	0.32 U
74-97-5	BROMOCHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
75-27-4	BROMODICHLOROMETHANE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U
75-25-2	BROMOFORM	ug/l	0.47 U	0.47 U	0.47 U	0.47 U
74-83-9	BROMOMETHANE	ug/l	0.2 UJ	0.2 UJ	0.2 UJ	0.2 UJ
75-15-0	CARBON DISULFIDE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
108-90-7	CHLOROBENZENE	ug/l	0.49 U	0.49 U	0.49 U	0.49 U
75-00-3	CHLOROETHANE	ug/l	0.2 U	0.2 UJ	0.2 U	0.2 U
67-66-3	CHLOROFORM	ug/l	0.34 U	0.34 U	0.34 U	0.34 U
74-87-3	CHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U
110-82-7	CYCLOHEXANE	ug/l	0.2 UJ	0.2 U	0.2 UJ	0.2 UJ
124-48-1	DIBROMOCHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
100-41-4	ETHYLBENZENE	ug/l	0.2 U	12.1	90	0.2 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/l	0.45 U	1.8 J	9.8	0.45 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/l	0.95 U	7.7 J	180	0.95 U
79-20-9	METHYL ACETATE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/l	1.3 U	29.2 J	1.3 U	1.3 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/l	2.1 U	200 J	2.1 U	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/l	0.2 U	0.2 UJ	0.44 J	0.2 U
75-09-2	METHYLENE CHLORIDE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/l	0.43 U	8.1	130	0.43 U
100-42-5	STYRENE	ug/l	0.36 U	0.49 J	0.36 U	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/l	0.35 U	0.35 U	1.6 J	1.6 J
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/l	0.27 U	0.27 U	0.27 U	0.27 U
108-88-3	TOLUENE	ug/l	0.37 U	36.9	62.8	0.37 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/l	0.29 U	0.29 U	0.29 U	0.29 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/l	0.28 U	0.28 U	0.28 U	0.28 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U
75-01-4	VINYL CHLORIDE	ug/l	0.34 U	0.34 U	0.34 U	0.34 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/l	100 U	100 U	100 U	100 U

G 1:1 / 1E	1.	Ir .: TD	MW 4	MW 5	MW 7	MW
Consolidated E		Location ID:	MW-4	MW-5	MW-7	MW-9
Ludlow Street S		Sample ID:	MW-4-20130408	MW-5-20130405	MW-7-20130404	MW-9-20130405
	ndwater Analytical Data	Lab Sample Id:	E1768-06	E1768-03	E1768-12	E1768-01
SDG: E1768		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	E1768	E1768	E1768	E1768
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	4/8/2013 8:40	4/5/2013 14:10	4/4/2013 12:32	4/5/2013 10:35
		Validated:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES		0.4.**			
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/l	0.4 U	0.253 U	0.206 U	0.21 U
	2,3,4,6-TETRACHLOROPHENOL	ug/l	0.4 U	0.253 U	0.206 U	0.21 U
	2,4,5-TRICHLOROPHENOL	ug/l	0.8 U	0.506 U	0.412 U	0.43 U
	2,4,6-TRICHLOROPHENOL	ug/l	1.1 U	0.709 U	0.577 U	0.6 U
	2,4-DICHLOROPHENOL	ug/l	1.3 U	0.835 U	0.68 U	0.7 U
	2,4-DIMETHYLPHENOL	ug/l	1.4 U	0.899 U	0.732 U	0.76 U
	2,4-DINITROPHENOL	ug/l	4.2 UJ	2.7 UJ	2.2 UJ	2.2 UJ
	2,4-DINITROTOLUENE	ug/l	2.1 U	1.3 U	1.1 U	1.1 U
606-20-2	2,6-DINITROTOLUENE	ug/l	0.64 U	0.405 U	0.33 U	0.34 U
91-58-7	2-CHLORONAPHTHALENE	ug/l	0.32 U	0.203 U	0.165 U	0.17 U
95-57-8	2-CHLOROPHENOL	ug/l	1.1 U	0.684 U	0.557 U	0.57 U
91-57-6	2-METHYLNAPHTHALENE	ug/l	0.64 U	42.1	340	0.34 U
	2-METHYLPHENOL (O-CRESOL)	ug/l	0.48 U	0.304 U	0.247 U	0.26 U
	2-NITROANILINE	ug/l	0.98 U	0.62 U	0.505 U	0.52 U
	2-NITROPHENOL	ug/l	1 U	0.658 U	0.536 U	0.55 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/l	4 U	R	2.1 U	2.1 U
МЕРНЗМЕРН	3+4-Methylphenols	ug/l	0.76 U	0.481 U	0.392 U	0.4 U
	3-NITROANILINE	ug/l	2.2 U	R	1.1 U	1.2 U
	4,6-DINITRO-2-METHYLPHENOL	ug/l	1.5 UJ	0.937 UJ	0.763 UJ	0.79 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/l	0.46 U	0.291 U	0.237 U	0.24 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/l	0.8 U	0.506 U	0.412 U	0.43 U
106-47-8	4-CHLOROANILINE	ug/l	5.7 U	3.6 U	2.9 U	3 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/l	0.42 U	0.266 U	0.216 U	0.22 U
	4-NITROANILINE	ug/l	2.7 U	1.7 UJ	1.4 U	1.4 U
100-01-0	4-NITROPHENOL	-	4 U	2.5 U	2.1 U	2.1 U
83-32-9	ACENAPHTHENE	ug/l	0.42 U	42.6	47.1	0.22 U
208-96-8		ug/l	0.42 U 1.4 U		38.6	0.22 U 0.74 U
	ACETORIENONE	ug/l		18.1		
98-86-2	ACETOPHENONE	ug/l	0.28 U	0.177 U	0.144 UJ	0.15 U
	ANTHRACENE	ug/l	0.32 U	0.203 U	26.2	0.17 U
	ATRAZINE	ug/l	0.8 U	0.506 U	0.412 U	0.43 U
	BENZALDEHYDE	ug/l	1.5 U	0.975 U	0.794 U	0.82 U
	BENZO(A)ANTHRACENE	ug/l	0.32 U	0.203 U	14.8	0.17 U
	BENZO(A)PYRENE	ug/l	0.28 U	0.177 U	9.7 J	0.15 U
	BENZO(B)FLUORANTHENE	ug/l	0.58 U	0.367 U	8.2 J	0.31 U
	BENZO(G,H,I)PERYLENE	ug/l	0.58 U	0.367 UJ	0.299 U	0.31 UJ
	BENZO(K)FLUORANTHENE	ug/l	0.36 U	0.228 U	0.186 U	0.19 U
	BENZYL BUTYL PHTHALATE	ug/l	0.38 U	0.241 U	0.196 U	0.2 U
	BIPHENYL (DIPHENYL)	ug/l	0.3 U	19.6	25.7	0.16 U
	BIS(2-CHLOROETHOXY) METHANE	ug/l	1.1 U	0.696 U	0.567 U	0.59 U
	BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/l	1.1 U	0.696 U	0.567 U	0.59 U
	BIS(2-CHLOROISOPROPYL) ETHER	ug/l	0.34 U	0.215 U	0.175 U	0.18 U
	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	0.32 U	0.203 U	0.165 U	0.17 U
	CAPROLACTAM	ug/l	4 U	2.5 U	2.1 U	2.1 U
86-74-8	CARBAZOLE	ug/l	0.44 U	0.278 U	5.5 J	0.23 U
218-01-9	CHRYSENE	ug/l	0.36 U	0.228 U	14.6	0.19 U
	DIBENZ(A,H)ANTHRACENE	ug/l	0.84 U	0.532 UJ	0.433 U	0.45 UJ
132-64-9	DIBENZOFURAN	ug/l	0.48 U	0.304 U	6.3 J	0.26 U
84-66-2	DIETHYL PHTHALATE	ug/l	0.76 U	0.481 U	0.392 U	0.4 U
	DIMETHYL PHTHALATE	ug/l	0.44 U	0.278 U	0.227 U	0.23 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	4 U	2.5 U	2.1 U	2.1 U
117-84-0	DI-N-OCTYLPHTHALATE	ug/l	1 U	0.646 U	0.526 U	0.54 U
206-44-0	FLUORANTHENE	ug/l	0.8 U	0.506 U	18.8	0.43 U
86-73-7	FLUORENE	ug/l	0.62 U	18	46.3	0.33 U
	HEXACHLOROBENZENE	ug/l	0.36 U	0.228 U	0.186 U	0.19 U
	HEXACHLOROBUTADIENE	ug/l	0.5 U	0.316 U	0.258 U	0.27 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/l	0.48 U	0.304 UJ	0.247 U	0.26 UJ
	HEXACHLOROETHANE	ug/l	0.5 U	0.316 U	0.258 U	0.27 U
	INDENO(1,2,3-C,D)PYRENE	ug/l	0.3 U	0.19 UJ	0.155 U	0.16 UJ
	ISOPHORONE	ug/l	0.6 U	0.38 U	0.309 U	0.32 U
	NAPHTHALENE	ug/l	0.24 U	240 J	1600	0.13 U
	NITROBENZENE	ug/l	1.4 U	0.861 U	0.701 U	0.72 U
	N-NITROSODI-N-PROPYLAMINE	ug/l	0.4 U	0.253 U	0.206 U	0.21 U
	N-NITROSODIPHENYLAMINE	ug/l	1.2 U	0.759 U	0.619 U	0.64 U
87-86-5	PENTACHLOROPHENOL	ug/l ug/l	3.4 UJ	2.2 UJ	1.8 UJ	1.8 UJ
	PHENANTHRENE	ug/l	0.52 U	21.2	110	0.28 U
	PHENOL	ug/l	0.32 U 0.42 U	0.266 U	0.216 U	0.28 U 0.22 U
	PYRENE	ug/l	0.4 U	0.253 U	37	0.21 U
127-00-0	LINENE	ug/I	0.4 0	U.433 U	31	0.21 0

C1: J-4- J	Edi	IIt' ID.	MW-4	MW 5	MW 7	MWO
Consolidated		Location ID:		MW-5	MW-7	MW-9
Ludlow Street		Sample ID:	MW-4-20130408	MW-5-20130405	MW-7-20130404	MW-9-20130405
	undwater Analytical Data	Lab Sample Id:	E1768-06	E1768-03	E1768-12	E1768-01
SDG: E1768		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	E1768	E1768	E1768	E1768
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	4/8/2013 8:40	4/5/2013 14:10	4/4/2013 12:32	4/5/2013 10:35
		Validated:	4/30/2013	4/30/2013	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:				
	TOTAL METALS					
7429-90-5	ALUMINUM	ug/l	32100 J	550 J	244 J	91.9 J
7440-36-0	ANTIMONY	ug/l	8 U	12.1 J	8 U	8 U
7440-38-2	ARSENIC	ug/l	40	4.4 J	4.2 U	9.05 J
7440-39-3	BARIUM	ug/l	523	916	192	380
7440-41-7	BERYLLIUM	ug/l	1.23 J	0.7 U	0.7 U	0.7 U
7440-43-9	CADMIUM	ug/l	1.3 J	2.2 J	R	R
7440-70-2	CALCIUM	ug/l	116000 J	267000 J	123000 J	196000 J
7440-47-3	CHROMIUM, TOTAL	ug/l	113 J	1.1 UJ	7.73 J	1.1 UJ
7440-47-3	COBALT	ug/l	25.1 J	5.8 UJ	5.8 UJ	5.8 UJ
7440-48-4	COPPER	ug/l	205 J	5.86 J	2.06 J	2 UJ
7439-89-6	IRON	ug/l	55300	4550	1300	11300
7439-89-6	LEAD		1200 J	232 J	2.6 UJ	2.6 UJ
7439-92-1	MAGNESIUM	ug/l	24500 J	59800 J	42200 J	77600 J
7439-95-4		ug/l				
	MANGANESE	ug/l	1210 J	1580 J	688 J	3240 J
7439-97-6	MERCURY	ug/l	20.182	0.092 U	0.092 U	0.092 U
7440-02-0	NICKEL	ug/l	112 J	4.2 UJ	9 J	4.2 UJ
7440-09-7	POTASSIUM	ug/l	11800	93500	26900	15100
7782-49-2	SELENIUM	ug/l	16.1	4.8 U	4.8 U	4.8 U
7440-22-4	SILVER	ug/l	1.5 UJ	1.5 UJ	1.5 UJ	1.5 UJ
7440-23-5	SODIUM	ug/l	712000 J	32700000 J	2230000 J	1210000 J
7440-28-0	THALLIUM	ug/l	4.02 J	2.4 UJ	2.4 UJ	3.46 J
7440-62-2	VANADIUM	ug/l	77.8 J	6.1 UJ	6.1 UJ	6.1 UJ
7440-66-6	ZINC	ug/l	715 J	46.8 J	10.3 J	6.5 UJ
	DISSOLVED METALS	1				
7429-90-5	ALUMINUM	ug/l	17.1 J			
7440-36-0	ANTIMONY	ug/l	9.22 J			
7440-38-2	ARSENIC	ug/l	6 J			
7440-39-3	BARIUM	ug/l	243			
7440-41-7	BERYLLIUM	ug/l	0.7 U			
7440-43-9	CADMIUM	ug/l	R			
7440-70-2	CALCIUM	ug/l	143000 J			
7440-47-3	CHROMIUM, TOTAL	ug/l	7.84 J			
7440-48-4	COBALT	ug/l	5.8 UJ			
7440-50-8	COPPER	ug/l	55.2 J			
7439-89-6	IRON	ug/l	60.5			
7439-92-1	LEAD	ug/l	3.29 J			
7439-95-4	MAGNESIUM	ug/l	21200 J			
7439-96-5	MANGANESE	ug/l	488 J			
7439-97-6	MERCURY	ug/l	0.092 U			
7440-02-0	NICKEL	ug/l	10.2 J			
7440-09-7	POTASSIUM	ug/l	8920			
7782-49-2	SELENIUM	ug/l	16.2			
7440-22-4	SILVER	ug/l	1.5 UJ			
7440-23-5	SODIUM	ug/l	753000 J			
7440-28-0	THALLIUM	ug/l	2.4 UJ			
7440-28-0	VANADIUM	ug/l	6.1 UJ			
7440-62-2	ZINC	ug/l	37.6 J			
7-7-00-0	OTHER	ug/1	37.03			
57-12-5	CYANIDE	ug/l	126	174	145	3 Ј
J 1 14"J	101111111111111111111111111111111111111	ug/1	120	1/7	177	ال ال

Consolidated I	Edison	Location ID:	FB040813	TRIPBLANK
Ludlow Street	Site	Sample ID:	FB040813-20130408	TBLANK-20130404
Validated Gro	undwater Analytical Data	Lab Sample Id:	E1768-10	E1768-11
SDG: E1768	•	Source:	CTECH	CTECH
		SDG:	E1768	E1768
		Matrix:	WATER	WATER
		Sampled:	4/8/2013 10:00	4/4/2013 9:00
		Validated:	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:		
	VOLATILES			
71-55-6	1,1,1-TRICHLOROETHANE	ug/l	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/l	0.31 U	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/l	0.45 U	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/l	0.38 U	0.38 U
75-34-3	1,1-DICHLOROETHANE	ug/l	0.36 U	0.36 U
75-35-4	1,1-DICHLOROETHENE	ug/l	0.47 U	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/l	0.2 U	0.2 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/l	0.2 U	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/l	0.46 U	0.46 U
106-93-4	1,2-DIBROMOETHANE (ETHYLENE DIBROMIDE)	ug/l	0.41 U	0.41 U
95-50-1 107-06-2	1,2-DICHLOROBENZENE 1,2-DICHLOROETHANE	ug/l	0.45 U 0.48 U	0.45 U 0.48 U
78-87-5	1,2-DICHLOROPROPANE	ug/l ug/l	0.48 U 0.46 U	0.48 U 0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/l ug/l	0.43 U	0.43 U
106-46-7	1.4-DICHLOROBENZENE	ug/l ug/l	0.43 U	0.43 U 0.32 U
591-78-6	2-HEXANONE	ug/l	1.9 U	1.9 U
67-64-1	ACETONE	ug/l	0.5 UJ	0.5 UJ
71-43-2	BENZENE	ug/l	0.3 U	0.32 U
71-43-2 74-97-5	BROMOCHLOROMETHANE	ug/l	0.32 U	0.32 U
75-27-4	BROMODICHLOROMETHANE	ug/l	0.2 U	0.36 U
75-25-2	BROMOFORM	ug/l	0.47 U	0.47 U
74-83-9	BROMOMETHANE	ug/l	0.2 UJ	0.2 UJ
75-15-0	CARBON DISULFIDE	ug/l	5.4	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/l	0.2 U	0.2 U
108-90-7	CHLOROBENZENE	ug/l	0.49 U	0.49 U
75-00-3	CHLOROETHANE	ug/l	0.2 U	0.2 U
67-66-3	CHLOROFORM	ug/l	0.34 U	0.34 U
74-87-3	CHLOROMETHANE	ug/l	0.2 U	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/l	0.35 U	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/l	0.31 U	0.31 U
110-82-7	CYCLOHEXANE	ug/l	0.2 UJ	0.2 UJ
124-48-1	DIBROMOCHLOROMETHANE	ug/l	0.2 U	0.2 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/l	0.2 U	0.2 U
100-41-4	ETHYLBENZENE	ug/l	0.2 U	0.2 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/l	0.45 U	0.45 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/l	0.95 U	0.95 U
79-20-9	METHYL ACETATE	ug/l	0.2 U	0.2 U
78-93-3	METHYL ETHYL KETONE (2-BUTANONE)	ug/l	1.3 U	1.3 U
108-10-1	METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	ug/l	2.1 U	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/l	0.2 U	0.2 U
75-09-2	METHYLENE CHLORIDE	ug/l	0.41 U	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/l	0.43 U	0.43 U
100-42-5	STYRENE	ug/l	0.36 U	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/l	0.35 U	0.35 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/l	0.27 U	0.27 U
108-88-3	TOLUENE TRANS 12 DICHI OPOETHENE	ug/l	0.37 U	0.37 U
156-60-5 10061-02-6	TRANS-1,2-DICHLOROETHENE TRANS-1,3-DICHLOROPROPENE	ug/l	0.41 U 0.29 U	0.41 U
		ug/l		0.29 U
79-01-6 75-60-4	TRICHLOROETHYLENE (TCE)	ug/l	0.28 U	0.28 U
75-69-4 75-01-4	TRICHLOROFLUOROMETHANE VINYL CHLORIDE	ug/l	0.35 U 0.34 U	0.35 U
		ug/l	U.34 U	0.34 U

Consolidated E		Location ID:	FB040813	TRIPBLANK
Ludlow Street S		Sample ID:	FB040813-20130408	TBLANK-20130404
	ndwater Analytical Data	Lab Sample Id:	E1768-10	E1768-11
SDG: E1768		Source:	CTECH	CTECH
		SDG:	E1768	E1768
		Matrix: Sampled:	WATER 4/8/2013 10:00	WATER 4/4/2013 9:00
		Validated:	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:	1/30/2013	1/30/2013
	SEMIVOLATILES			
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/l	0.217 U	
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/l	0.217 U	
	2,4,5-TRICHLOROPHENOL	ug/l	0.435 U	
	2,4,6-TRICHLOROPHENOL	ug/l	0.609 U	
	2,4-DICHLOROPHENOL 2,4-DIMETHYLPHENOL	ug/l ug/l	0.717 U 0.772 U	
	2,4-DINITROPHENOL	ug/l	2.3 UJ	
	2,4-DINITROTOLUENE	ug/l	1.1 U	
	2,6-DINITROTOLUENE	ug/l	0.348 U	
91-58-7	2-CHLORONAPHTHALENE	ug/l	0.174 U	
95-57-8	2-CHLOROPHENOL	ug/l	0.587 U	
91-57-6	2-METHYLNAPHTHALENE	ug/l	0.348 U	
	2-METHYLPHENOL (O-CRESOL)	ug/l	0.261 U	
	2-NITROANILINE	ug/l	0.533 U	
88-75-5	2-NITROPHENOL	ug/l	0.565 U	
	3,3'-DICHLOROBENZIDINE	ug/l	2.2 U	
MEPH3MEPH 99-09-2	3+4-Methylphenols 3-NITROANILINE	ug/l ug/l	0.413 U 1.2 U	
534-52-1	4.6-DINITRO-2-METHYLPHENOL	ug/l	0.804 UJ	
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/l	0.25 U	
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/l	0.435 U	
106-47-8	4-CHLOROANILINE	ug/l	3.1 U	
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/l	0.228 U	
100-01-6	4-NITROANILINE	ug/l	1.5 U	
100-02-7	4-NITROPHENOL	ug/l	2.2 U	
83-32-9	ACENAPHTHENE	ug/l	0.228 U	
208-96-8	ACENAPHTHYLENE	ug/l	0.761 U	
98-86-2	ACETOPHENONE	ug/l	0.152 U	
120-12-7 1912-24-9	ANTHRACENE ATRAZINE	ug/l ug/l	0.174 U 0.435 U	
	BENZALDEHYDE	ug/l	0.433 U 0.837 U	
	BENZO(A)ANTHRACENE	ug/l	0.174 U	
	BENZO(A)PYRENE	ug/l	0.152 U	
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.315 U	
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.315 U	
	BENZO(K)FLUORANTHENE	ug/l	0.196 U	
	BENZYL BUTYL PHTHALATE	ug/l	0.207 U	
	BIPHENYL (DIPHENYL)	ug/l	0.163 U	
	BIS(2-CHLOROETHOXY) METHANE BIS(2-CHLOROETHYL) ETHER (2-CHLOROETHYL ETHER)	ug/l	0.598 U	
	BIS(2-CHLOROISOPROPYL) ETHER BIS(2-CHLOROISOPROPYL) ETHER	ug/l ug/l	0.598 U 0.185 U	
	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	0.174 U	
105-60-2	CAPROLACTAM	ug/l	2.2 U	
86-74-8	CARBAZOLE	ug/l	0.239 U	
218-01-9	CHRYSENE	ug/l	0.196 U	
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.457 U	
	DIBENZOFURAN	ug/l	0.261 U	
	DIETHYL PHTHALATE	ug/l	0.413 U	
	DIMETHYL PHTHALATE	ug/l	0.239 U	
	DI-N-BUTYL PHTHALATE	ug/l	2.2 U	
	DI-N-OCTYLPHTHALATE	ug/l	0.554 U	
	FLUORANTHENE FLUORENE	ug/l	0.435 U 0.337 U	
	HEXACHLOROBENZENE	ug/l ug/l	0.196 U	
	HEXACHLOROBUTADIENE	ug/l	0.272 U	
	HEXACHLOROCYCLOPENTADIENE	ug/l	0.261 U	
	HEXACHLOROETHANE	ug/l	0.272 U	
	INDENO(1,2,3-C,D)PYRENE	ug/l	0.163 U	
	ISOPHORONE	ug/l	0.326 U	
	NAPHTHALENE	ug/l	0.13 U	
	NITROBENZENE	ug/l	0.739 U	
	N-NITROSODI-N-PROPYLAMINE	ug/l	0.217 U	
	N-NITROSODIPHENYLAMINE	ug/l	0.652 U	
	PENTACHLOROPHENOL DHENANTHRENE	ug/l	1.9 UJ	
	PHENANTHRENE	ug/l	0.283 U	
	PHENOL	ug/l	0.228 U	

C1: 1-4-1	F.E	IIti ID.	ED040012	TRIPRI ANIZ
Consolidated		Location ID:	FB040813	TRIPBLANK
Ludlow Stree		Sample ID:	FB040813-20130408	TBLANK-20130404
	oundwater Analytical Data	Lab Sample Id:	E1768-10	E1768-11
SDG: E1768		Source:	CTECH	CTECH
		SDG:	E1768	E1768
		Matrix:	WATER	WATER
		Sampled:	4/8/2013 10:00	4/4/2013 9:00
		Validated:	4/30/2013	4/30/2013
CAS NO.	COMPOUND	UNITS:		
	TOTAL METALS			
7429-90-5	ALUMINUM	ug/l	6.5 UJ	
7440-36-0	ANTIMONY	ug/l	8 U	
7440-38-2	ARSENIC	ug/l	4.2 U	
7440-39-3	BARIUM	ug/l	4 U	
7440-41-7	BERYLLIUM	ug/l	0.7 U	
7440-43-9	CADMIUM	ug/l	R	
7440-70-2	CALCIUM	ug/l	99.7 J	
7440-47-3	CHROMIUM, TOTAL	ug/l	1.1 UJ	
7440-48-4	COBALT	ug/l	5.8 UJ	
7440-50-8	COPPER	ug/l	2 UJ	
7439-89-6	IRON	ug/l	27.8 J	
7439-92-1	LEAD	ug/l	2.6 UJ	
7439-95-4	MAGNESIUM	ug/l	35.3 J	
7439-96-5	MANGANESE	ug/l	1.7 U	
7439-97-6	MERCURY	ug/l	0.092 U	
7440-02-0	NICKEL	ug/l	4.2 UJ	
7440-09-7	POTASSIUM	ug/l	594 J	
7782-49-2	SELENIUM	ug/l	4.8 U	
7440-22-4	SILVER	ug/l	1.5 UJ	
7440-23-5	SODIUM	ug/l	34500 J	
7440-28-0	THALLIUM	ug/l	2.4 UJ	
7440-62-2	VANADIUM	ug/l	6.1 UJ	
7440-66-6	ZINC	ug/l	6.5 UJ	
	DISSOLVED METALS			
7429-90-5	ALUMINUM	ug/l	6.5 UJ	
7440-36-0	ANTIMONY	ug/l	8 U	
7440-38-2	ARSENIC	ug/l	4.2 U	
7440-39-3	BARIUM	ug/l	4 U	
7440-41-7	BERYLLIUM	ug/l	0.7 U	
7440-43-9	CADMIUM	ug/l	R	
7440-70-2	CALCIUM	ug/l	210 J	
7440-47-3	CHROMIUM, TOTAL	ug/l	1.1 UJ	
7440-48-4	COBALT	ug/l	5.8 UJ	
7440-50-8	COPPER	ug/l	2 UJ	
7439-89-6	IRON	ug/l	20.4 U	
7439-92-1	LEAD	ug/l	2.6 UJ	
7439-95-4	MAGNESIUM	ug/l	32.5 U	
7439-95-4	MANGANESE	ug/l	2.59 J	
7439-97-6	MERCURY	ug/l	0.092 U	
7440-02-0	NICKEL	ug/l	4.2 UJ	
7440-02-0	POTASSIUM	ug/l ug/l	273 J	
7782-49-2	SELENIUM	ug/l	4.8 U	
7440-22-4	SILVER		1.5 UJ	
7440-22-4	SODIUM	ug/l	1.5 UJ 3160 J	
		ug/l		
7440-28-0	THALLIUM	ug/l	2.4 UJ	
7440-62-2	VANADIUM	ug/l	6.1 UJ	
7440-66-6	ZINC OTHER	ug/l	26.9 J	
57-12-5	CYANIDE	ug/l	3 U	
01-12-0	CITITULE	ug/1	3.0	l